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STRUCTURE FILE UPDATES: 12 APR 2012 HIGHEST RN 1367421-69-9 DICTIONARY FILE UPDATES: 12 APR 2012 HIGHEST RN 1367421-69-9

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TSCA INFORMATION NOW CURRENT THROUGH DECEMBER 23, 2011

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http://www.cas.org/support/stngen/stndoc/properties.html

=> d que 138

L5 STR

SO3H @6 G17 CH2=C-1CH2-O

CH2=C-15 012 13 14 G2 012 13 14 16

G3 17

VAR G1=1/6
VAR G2=O/N/S
VAR G3=8/12
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DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 17

STEREO ATTRIBUTES: NONE

L7 28963 SEA FILE=REGISTRY SSS FUL L5

L10 STR

CH2: C - CH2-0 - G12 7 7 A @13

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GRAPH ATTRIBUTES:

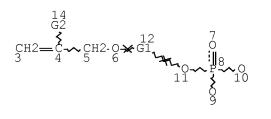
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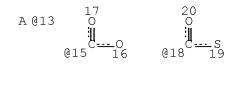
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L12 797 SEA FILE=REGISTRY SUB=L7 SSS FUL L10

L33 STF





23 0 :::: C----N @21 22

REP G1=(1-20) 13

VAR G2=15/18/21/COOH/SO3H

NODE ATTRIBUTES:

NSPEC IS RC AT 13 DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 21

STEREO ATTRIBUTES: NONE

L35 24 SEA FILE=REGISTRY SUB=L12 SSS FUL L33

L37 8 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L35

L38 3 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L37 AND (1802-2003

)/PRY,AY,PY

=> d 138 1-3 ibib ed abs hitstr hitind
YOU HAVE REQUESTED DATA FROM FILE 'HCAPLUS' - CONTINUE? (Y)/N:n

=> fil hcap FILE 'HCAPLUS' ENTERED AT 11:19:19 ON 13 APR 2012 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2012 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 13 Apr 2012 VOL 156 ISS 17

FILE LAST UPDATED: 12 Apr 2012 (20120412/ED)

REVISED CLASS FIELDS (/NCL) LAST RELOADED: Dec 2011

USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Dec 2011

HCAplus now includes complete International Patent Classification (IPC) reclassification data for the fourth quarter of 2011.

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> d 138 1-3 ibib ed abs hitstr hitind

L38 ANSWER 1 OF 3 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2005:568976 HCAPLUS Full-text

DOCUMENT NUMBER: 143:83603

TITLE: One-part self-etching, self-priming dental

adhesive composition

INVENTOR(S): Klee, Joachim E.; Lehmann, Uwe; Walz, Uwe

PATENT ASSIGNEE(S): Dentsply Detrey GmbH, Germany

SOURCE: Eur. Pat. Appl., 30 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

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             KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW,
             MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD,
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             DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC,
             NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA,
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PRIORITY APPLN. INFO.:
                                             EP 2003-29824
                                                                 Α
                                                                   20031223
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                                             WO 2004-EP14307
                                                                    20041215
                                                                 M
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ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

ED Entered STN: 01 Jul 2005

One-part self-etching, self-priming dental adhesive compn. having a pH of at most 2 comprises (a) a polymerizable acidic phosphoric acid ester monomer; (b) one or more polymerizable acidic monomers; (c) a polymerizable N-substituted alkylacrylic or acrylic acid amide monomer; (d) an org. and/or inorg. acid; (e) an org. water sol. solvent and/or water; and (f) polymn. initiator, inhibitor and stabilizer. An adhesive polymer was prepd. from 2-acrylamido-2-methyl-propane-sulfonic acid, 3, (4), 8, (9)-bis(acrylamido methyl) tricyclo-5.2.1.02,6 decane, Et 2-[13-dihydrogen phosphoryl-13,2-dioxatridecyl]acrylate, and N,N'-bisacrylamido-N,N'-diethyl-1,3-propane.

IT 752234-98-3P 752235-00-0P 855894-56-3P

(one-part self-etching, self-priming dental adhesive compn.)

RN 752234-98-3 HCAPLUS

CN 2-Propenoic acid, 2-[[[10-(phosphonooxy)decyl]oxy]methyl]-, 1-ethyl ester (CA INDEX NAME)

RN 752235-00-0 HCAPLUS

CN 2-Propenoic acid, 2-[[2-(phosphonooxy)ethoxy]methyl]-, 1-ethyl ester (CA INDEX NAME)

RN 855894-56-3 HCAPLUS

CN 2-Propenoic acid, 2-[[2-(phosphonooxy)ethoxy]methyl]- (CA INDEX NAME)

$$\begin{array}{c} & \text{CH2} \\ \text{HO2C-C-CH2-O-CH2-CH2-OPO3H2} \end{array}$$

IT 855894~57~4P, 2-Acrylamido-2-methyl-propane-sulfonic acid-3,(4),8,(9)-bis(acrylamido methyl) tricyclo-5.2.1.02,6 decane-Ethyl 2-[13-dihydrogen phosphoryl-13,2-dioxatridecyl]acrylate-N,N'-Bisacrylamido-N,N'-diethyl-1,3-propane copolymer 855894~58~5P

(one-part self-etching, self-priming dental adhesive compn.)

RN 855894-57-4 HCAPLUS

CN 2-Propenoic acid, 2-[[[10-(phosphonooxy)decyl]oxy]methyl]-, 1-ethyl ester, polymer with 2-methyl-2-[(1-oxo-2-propen-1-yl)amino]-1-propanesulfonic acid, N,N-[[octahydro-4,7-methano-1H-indene-1,5(1,6 or 2,5)diyl]bis(methylene)]bis[2-propenamide] and N,N'-1,3-propanediylbis[N-ethyl-2-propenamide] (CA INDEX NAME)

CM 1

CRN 855532-00-2 CMF C18 H26 N2 O2 CCI IDS

CM 2

CRN 752234-98-3 CMF C16 H31 O7 P

CM 3

CRN 442200-41-1 CMF C13 H22 N2 O2

CM 4

CRN 15214-89-8 CMF C7 H13 N O4 S

RN 855894-58-5 HCAPLUS

CN 2-Propenoic acid, 2-[[2-(phosphonooxy)ethoxy]methyl]-, 1-ethyl ester, polymer with 2-methyl-2-[(1-oxo-2-propen-1-yl)amino]-1-propanesulfonic acid, N,N-[[octahydro-4,7-methano-1H-indene-1,5(1,6 or 2,5)diyl]bis(methylene)]bis[2-propenamide] and N,N'-1,3-propanediylbis[N-ethyl-2-propenamide] (CA INDEX NAME)

CM 1

CRN 855532-00-2 CMF C18 H26 N2 O2 CCI IDS

CM 2

CRN 752235-00-0 CMF C8 H15 O7 P

CM 3

CRN 442200-41-1 CMF C13 H22 N2 O2

CM 4

CRN 15214-89-8 CMF C7 H13 N O4 S

IPCI C07F0009-00 [I,C]; A61K0006-00 [I,C]; A61K0006-02 [I,C]; C08F0030-00

[I,C]; C07F0009-09 [I,A]; A61K0006-00 [I,A]; A61K0006-083 [I,A]; C08F0030-02 [I,A] IPCR A61K0006-00 [I,A]; A61K0006-083 [I,A]; C07F0009-09 [I,A]; C08F0030-02 [I,A]63-8 (Pharmaceuticals) CC 752234-97-2P **752234-98-3P** 752234-99-4P ΙT 752235-00-0P 855894-56-3P (one-part self-etching, self-priming dental adhesive compn.) 855894-57-4P, 2-Acrylamido-2-methyl-propane-sulfonic ΙT acid-3, (4),8,(9)-bis(acrylamido methyl) tricyclo-5.2.1.02,6 decane-Ethyl 2-[13-dihydrogen phosphoryl-13,2-dioxatridecyl]acrylate-N, N'-Bisacrylamido-N, N'-diethyl-1, 3-propane copolymer 855894-58-5P (one-part self-etching, self-priming dental adhesive compn.) OS.CITING REF COUNT: 7 THERE ARE 7 CAPLUS RECORDS THAT CITE THIS RECORD (7 CITINGS) REFERENCE COUNT: THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT L38 ANSWER 2 OF 3 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2005:182218 HCAPLUS <u>Full-text</u> 142:287808 DOCUMENT NUMBER: Lithographic printing plate precursor for direct TITLE: imaging from a digital data and developing in a printing machine without passing through a development step INVENTOR(S): Yamasaki, Sumiaki; Makino, Naonori; Inno, Toshifumi Fuji Photo Film Co., Ltd., Japan PATENT ASSIGNEE(S): U.S. Pat. Appl. Publ., 50 pp. SOURCE: CODEN: USXXCO DOCUMENT TYPE: Patent English LANGUAGE: FAMILY ACC. NUM. COUNT: 1 PATENT INFORMATION: PATENT NO. APPLICATION NO. DATE KIND DATE ---------_____ US 20050048398 A1 20050303 US 2004-896070 20040722 <--B2 20070227 A2 20050126 EP 2004-17306 US 7183038 EP 1500498 20040722 <--A3 20051012 B1 20101215 EP 1500498 EP 1500498 R: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LI, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR JP 2005238816 Α 20050908 JP 2004-214190 20040722 <--AT 491968 Τ 20110115 AT 2004-17306 20040722 <--PRIORITY APPLN. INFO.: JP 2003-277448 A 20030722

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JP 2004-652 A 20040105

JP 2004-17599 A 20040126

JP 2004-214190 A 20040722

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 142:287808

ED Entered STN: 04 Mar 2005

GΙ

• Br-

AB A lithog, printing plate precursor is described for recording an image directly from a digital data and development in a printing machine without passing through a development step. The precursor provides lithog, printing plates with improved press life and stain resistance. Thus, the precursor coating compn. comprises an image-forming layer contg. a polymn. initiator and a polymerizable compd., and a hydrophilic support. The compn. includes a compd. contg. at least one functional group interacting with a surface of the hydrophilic support. This compd. is one of a phosphonic acid and a phosphoric acid amide.

II 847226-71-5

(lithog. printing plate precursor for direct imaging from digital data and in-press development)

RN 847226-71-5 HCAPLUS

CN 2-Propenoic acid, 2-(15,15-dihydroxy-15-oxido-2,5,8,11,14-pentaoxa-15-phosphapentadec-1-yl)-, 1-ethyl ester (CA INDEX NAME)

PAGE 1-B

-- OPO3H2

IT 847204-83-5 847204-84-6

(phosphonic deriv.; lithog. printing plate precursor for direct imaging from digital data and in-press development)

RN 847204-83-5 HCAPLUS

CN 2-Propenoic acid, 2-(9,9-dihydroxy-9-oxido-2,5,8-trioxa-9-phosphanon-1-yl)-, 1-methyl ester (CA INDEX NAME)

RN 847204-84-6 HCAPLUS

CN 2-Propenoic acid, 2-(15,15-dihydroxy-15-oxido-2,5,8,11,14-pentaoxa-15-phosphapentadec-1-yl)-, 1-methyl ester (CA INDEX NAME)

PAGE 1-A

PAGE 1-B

- OPO3H2

INCL 430270100

IPCI G03C0001-77 [I,A]; G03C0001-91 [I,A]; G03F0007-028 [I,A]; G03F0007-038 [I,A]; G03F0007-09 [I,A]

IPCR G03F0007-027 [I,A]; B41C0001-10 [I,A]; B41N0001-08 [I,A]; B41N0001-14
 [I,A]; B41N0003-03 [I,A]; G03F0007-00 [I,A]; G03C0001-77 [I,A];
 G03C0001-91 [I,A]; G03F0007-028 [I,A]; G03F0007-038 [I,A]; G03F0007-09
 [I,A]

NCL 430/270.100; 430/271.100; 101/456.000; 101/457.000; 101/459.000; 101/467.000; 430/278.100; 430/281.100; 430/302.000; 430/325.000; 430/944.000

CC 74-5 (Radiation Chemistry, Photochemistry, and Photographic and Other Reprographic Processes)

IT 847204-87-9 847204-88-0 847204-89-1 847204-90-4 847204-91-5 847226-71-5

(lithog. printing plate precursor for direct imaging from digital data and in-press development)

IT 80730-17-2 223681-84-3 847204-73-3 847204-74-4 847204-75-5 847204-76-6 847204-77-7 847204-78-8 847204-82-4 847204-83-5 847232-64-8

RECORD (8 CITINGS)

REFERENCE COUNT: 25 THERE ARE 25 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L38 ANSWER 3 OF 3 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2004:732258 HCAPLUS Full-text

DOCUMENT NUMBER: 141:243056

TITLE: Polymerizable phosphoric acid ester derivatives

for dental compositions

INVENTOR(S): Klee, Joachim E.; Lehmann, Uwe; Walz, Uwe; Liu,

Huaibing

PATENT ASSIGNEE(S): Dentsply Detrey GmbH, Germany

SOURCE: Eur. Pat. Appl., 20 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

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	WO	2004	0781	00		A2		2004	0916		WO 2		 EP22	89		2	0040305
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	WO	2004				A3		2004									
		W:	•	•	•	•	•	AU,	•	•	•	•	•	•	•	•	•
				-				CZ,	•	-	-		-		-	-	· •
				•				HR,	•	•	•		•		•	•	•
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		RW:	•	•	•	•	•	MW,	•	•	•	•	•	•	•	•	•
								DE,									
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		1.601	C70			D 1		2011	O F 1 1			<					
	EΡ	1601				B1		2011		C.D.	C.D.					ο. Ε	140
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	TD	2006	PL,					2006	0007		TD 0	000	- O 4 -	C 2		^	0040205
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PRIOR	ΙΤΊ	APP	LN.	INFO	.:						EP 2					A 2	0030307
											WO 2	004-	EP22	89		W 2	0040305

ED Entered STN: 09 Sep 2004

- AB The present invention provides a polymerizable phosphoric acid ester deriv. for use in dental compns. E.g., 2,2-tris(2,6-dioxa-4-methylene-5-oxo-octyl)ethanol phosphoric acid ester was
 - 2,2,2-tris(2,6-dioxa-4-methylene-5-oxo-octyl)ethanol phosphoric acid ester was prepd. from pentaerythritol, Et chloromethyacrylate, and then treatment with the product with POCl3 and hydrolyzed.
- IT 752234-96-1P 752234-98-3P 752235-00-0P

(polymerizable phosphoric acid ester derivs. for dental compns.)

- RN 752234-96-1 HCAPLUS
- CN 2-Propenoic acid, 2,2'-[[2-[[[2-(ethoxycarbonyl)-2-propenyl]oxy]methyl]-2-[(phosphonooxy)methyl]-1,3-propanediyl]bis(oxymethylene)]bis-, 1,1'-diethyl ester (9CI) (CA INDEX NAME)

- RN 752234-98-3 HCAPLUS
- CN 2-Propenoic acid, 2-[[[10-(phosphonooxy)decyl]oxy]methyl]-, 1-ethyl ester (CA INDEX NAME)

- RN 752235-00-0 HCAPLUS
- CN 2-Propenoic acid, 2-[[2-(phosphonooxy)ethoxy]methyl]-, 1-ethyl ester (CA INDEX NAME)

- IPCI C07F0009-09 [ICM,7]; A61K0006-08 [ICS,7]; C08F0030-02 [ICS,7]
- IPCR A61K0006-00 [I,A]; A61K0006-08 [I,A]; A61K0006-083 [I,A]; C07F0009-09
 [I,A]; C08F0030-02 [I,A]
- CC 23-17 (Aliphatic Compounds)
 - Section cross-reference(s): 63
- IT 752234-96-1P 752234-98-3P 752235-00-0P
- (polymerizable phosphoric acid ester derivs. for dental compns.)
- OS.CITING REF COUNT: 3 THERE ARE 3 CAPLUS RECORDS THAT CITE THIS
 - RECORD (3 CITINGS)
- REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR
 - THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

=> d que 139

L5 STR

SO3H@6 G17

CH2=C~CH2·O

CH2=C-C-G2 @12 13 14 16

G3 17

VAR G1=1/6 VAR G2=O/N/S VAR G3=8/12 NODE ATTRIBUTES: DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 17

STEREO ATTRIBUTES: NONE

L7 28963 SEA FILE=REGISTRY SSS FUL L5

L10 STR

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NSPEC IS RC AT 13
DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 11

STEREO ATTRIBUTES: NONE

L12 797 SEA FILE=REGISTRY SUB=L7 SSS FUL L10

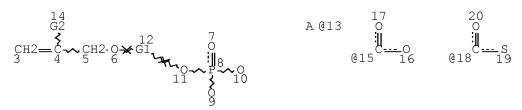
L16 STR

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GRAPH ATTRIBUTES:
RING(S) ARE ISOLATED OR EMBEDDED
NUMBER OF NODES IS 13

STEREO ATTRIBUTES: NONE

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L22	400	SEA FI	LE=HCAPLUS	SPE=ON	ABB=ON	PLU=ON	L12	
L23	13991	SEA FI	LE=HCAPLUS	SPE=ON	ABB=ON	PLU=ON	L18	
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GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 21

STEREO ATTRIBUTES: NONE

L35 24 SEA FILE=REGISTRY SUB=L12 SSS FUL L33

L37 8 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L35

L38 3 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L37 AND (1802-2003

)/PRY,AY,PY

L39 19 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L30 NOT L38

=> d 139 1-19 ibib ed abs hitstr hitind

L39 ANSWER 1 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2005:33579 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 142:115528

TITLE: Ethylene-vinyl acetate copolymer emulsions, their

manufacture, and adhesive compositions thereof with balanced heat resistance and low-temperature

properties

INVENTOR(S): Yako, Manabu; Yamamoto, Hiroki
PATENT ASSIGNEE(S): Denki Kagaku Kogyo Co., Ltd., Japan

SOURCE: Jpn. Kokai Tokkyo Koho, 13 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 2005008718	A	20050113	JP 2003-172880	20030618
JP 4041534 PRIORITY APPLN. INFO.:	В2	20080130	JP 2003-172880 <	20030618

OTHER SOURCE(S): MARPAT 142:115528

ED Entered STN: 14 Jan 2005

The emulsions have av. particle size .gtoreq.3 .mu.m and nonvolatile fraction .gtoreq.70% and are manufd. by using 0.2-3.0 parts/100 parts vinyl acetate (I) of reactive surfactants having structures H2C:CHCH2OCOCH(SO3Na)CH2CO2R (R = H, alkyl), H2C:CHCH2OCCH2CH(CH2OC6H4R) (OCH2CH2) nOX (n = 1-50; X = H, SO3NH4; R = C5-12 alkyl), or [H2C:CHCH2O(CH2CH2MeO)n]mP(O) (OH)3-m (n = 1-50; m = 1-2) and 0.5-4.0 parts/100 parts I of vinyl alc. polymer (PVA) with sapon. degree 60-85 mol% and av. d.p. 100-500 as emulsifiers in polymn. Thus, 70 parts I and 25 parts ethylene were polymd. in an aq. soln. contg. 3 parts PVA (UMR 10H; sapon. degree 80 mol%, d.p. 400), 1 part sulfosuccinate-type reactive emulsifier (Eleminol JS 2), nonionic surfactants (Emulgen 913, Emulgen 950), NaOAc, Rongalit, FeSO4.7H2O, EDTA.4Na, and ammonium persulfate, then 30 parts I was added dropwise to the soln. and polymd. to give an EVA emulsion showing nonvolatile fraction 71.1%, av. particle size 5.1 .mu.m, viscosity 1800 mPa-s at 30.degree. and 30 rpm, and PhMe-insol. fraction 42%.

IT 820213-85-2P

(manuf. of ethylene-vinyl acetate polymer emulsions by using reactive emulsifier and poly(vinyl alc.) for cold- and heat-resistant adhesives)

RN 820213-85-2 HCAPLUS

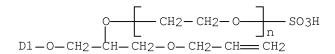
CN Acetic acid ethenyl ester, polymer with ethene and .alpha.-sulfo-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt, graft (9CI) (CA INDEX NAME)

CM 1

CRN 113405-85-9 CMF (C2 H4 O)n C21 H34 O6 S . H3 N CCI IDS, PMS



 $D1 - (CH_2) 8 - Me$



● NH3

CM 2

CRN 108-05-4 CMF C4 H6 O2

AcO-CH = CH₂

CM 3

CRN 74-85-1 CMF C2 H4

H2C==CH2

IT 86271-46-7, Allyl sodium sulfosuccinate 220308-52-1, Polypropylene glycol monoallyl ether, phosphate

(reactive emulsifier; manuf. of ethylene-vinyl acetate polymer emulsions by using reactive emulsifier and poly(vinyl alc.) for cold- and heat-resistant adhesives)

RN 86271-46-7 HCAPLUS

CN Butanedioic acid, 2-sulfo-, 1-(2-propen-1-yl) ester, sodium salt (1:1) (CA INDEX NAME)

Na

RN 220308-52-1 HCAPLUS

CN Poly[oxy(methyl-1,2-ethanediyl)],
.alpha.-phosphono-.omega.-(2-propen-1-yloxy)- (CA INDEX NAME)

$$H_2C$$
 = CH - CH_2 - O - O

IPCI C08F0210-00 [I,A]; C08F0218-08 [I,A]; C08F0002-24 [I,A]; C09J0131-04
 [I,A]; C09J0011-08 [I,A]

IPCR C08F0002-24 [I,A]; C08F0010-00 [I,A]; C09J0011-08 [I,A]; C09J0131-04
 [I,A]; C08F0210-00 [I,A]; C08F0218-08 [I,A]

CC 38-3 (Plastics Fabrication and Uses)

IT 820213-84-1P, Eleminol JS 2-ethylene-vinyl acetate copolymer 820213-85-2P 820233-84-9P

(manuf. of ethylene-vinyl acetate polymer emulsions by using reactive emulsifier and poly(vinyl alc.) for cold- and heat-resistant adhesives)

IT 86271-46-7, Allyl sodium sulfosuccinate

220308-52-1, Polypropylene glycol monoallyl ether, phosphate (reactive emulsifier; manuf. of ethylene-vinyl acetate polymer emulsions by using reactive emulsifier and poly(vinyl alc.) for cold- and heat-resistant adhesives)

L39 ANSWER 2 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2004:856797 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 141:350862

TITLE: Reactive liquid polymer crosslinking agent and

process for preparation

INVENTOR(S): Lazar, Warren G.; Clark, James A.

PATENT ASSIGNEE(S): USA

SOURCE: U.S. Pat. Appl. Publ., 18 pp., Cont.-in-part of

U.S. Ser. No. 13,164, abandoned.

CODEN: USXXCO

DOCUMENT TYPE: Patent LANGUAGE: English

10/596.747

FAMILY ACC. NUM. COUNT: 2 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 20040200993	A1	20041014	US 2004-833816	20040427
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US 20030168629	A1	20030911	US 2001-13164	20011210
			<	
PRIORITY APPLN. INFO.:			US 2001-13164	B2 20011210
			<	

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

ED Entered STN: 18 Oct 2004

AB A reactive liq. crosslinking agent for use in the prepn. of polymeric substances. The crosslinking agent comprises a substituted 1,3,5-triazine reacted with H2O, an acid alkyl sulfonate and/or phosphonate and a solidifying modifier contg. an hydroxyl functional group to form a substituted 1,3,5-triazine hydrate. The reactive liq. polymer crosslinking agent has a solids content between 20-99% solids. The reactive liq. crosslinking agents (RLPC's) are useful as modifiers in the prepn. of polymeric compds. which are suitable for 1-component self-crosslinking adhesives, coatings and polymers used in optics, textiles, composites, casting and molding. RLPC systems contg. from 1-30% RLPC provide fast single package thermosetting polymeric compds. with enhanced properties such as chem., heat and abrasion resistance.

RN 15214-89-8 HCAPLUS

CN 1-Propanesulfonic acid, 2-methyl-2-[(1-oxo-2-propen-1-yl)amino]- (CA INDEX NAME)

RN 111083-74-0 HCAPLUS

CN Titanate(3-), [2,2-bis[(2-propen-1-yloxy)methyl]-1-butanolato-.kappa.0][P,P-dioctyl diphosphato(2-)-.kappa.0'']bis[P,P-dioctyl diphosphato(2-)-.kappa.0'',.kappa.0''']-, hydrogen (1:3) (CA INDEX NAME)

PAGE 1-A

CH2-O-CH2-CH=CH2

PAGE 2-A

●з н+

INCL 252182130

IPCI C09K0003-00 [ICM,7]

IPCR C08F0008-30 [I,A]; C09D0167-00 [I,A]; C09J0167-00 [I,A]

NCL 252/182.130

37-6 (Plastics Manufacture and Processing) ΙT 57-50-1D, Sucrose, alkylglycosides, reaction products with triazine and sulfonylzirconate 98-11-3D, Phenylsulfonic acid, reaction products with triazine and diethylene glycol 107-21-1D, Ethylene glycol, reaction products with triazine and sulfonyltitanate 108-78-1D, 2,4,6-Triamino-1,3,5-triazine, reaction products with phenylphosphoric acid 110-63-4D, 1,4-Butanediol, reaction products with triazine and phosphatotitanate 111-46-6D, Diethylene glycol, reaction products with phenylsulfonic acid and triazine 504-63-2D, 1,3-Propylene glycol, reaction products with triazine and sulfonate 629-11-8D, 1,6-Hexanediol, reaction products with triazine and phosphatotitanate 1571-33-1D, Phenylphosphonic acid, reaction products with triazine 5606-17-7D, reaction products with sulfonate and propylene glycol 5606-19-9D, reaction products with polypropylene glycol and sulfate ester 15214-89-80, 2-Acrylamido-2-methylpropanesulfonic acid, reaction products with triazine and polyethylene glycol 25322-68-3D, Polyethylene glycol, reaction products with triazine and sulfonate 25322-69-4D, Polypropylene glycol, reaction products with triazine and sulfate 89619-91-0D, reaction products with alkylglycosides and ester sulfonylzirconate 103406-74-2D, reaction products with triazine and ethylene glycol 109766-35-0D, reaction products with triazine and alkylglycosides 111083-74-0D, reaction products with triazine and diol 544651-50-5D, reaction products with sulfonate and polyethylene glycol 544651-51-6D, reaction products with phosphate

ester and polyethylene glycol 544651-52-7D, reaction products with phosphotitanate and diol

(reactive liq. polymer crosslinking agent)

L39 ANSWER 3 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2004:467829 HCAPLUS Full-text

DOCUMENT NUMBER: 141:27260

TITLE: Defoaming agent for cementitious composition INVENTOR(S): Lorenz, Klaus; Yaguchi, Minoru; Sugiyama, Tomomi;

Albrecht, Gerhard

PATENT ASSIGNEE(S): Die Construction Research & Technology G.m.b.H.,

Germany

SOURCE: PCT Int. Appl., 22 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.																	
	WO 2004048293				A1 20040610							0021					
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		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DΖ,	EC,	EE,	ES,	FΙ,	GB,	GD,	
		GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KΕ,	KG,	KΡ,	KR,	KΖ,	
		LC,	LK,	LR,	LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	
		NO,	NΖ,	OM,	PH,	PL,	PT,	RO,	RU,	SD,	SE,	SG,	SI,	SK,	SL,	ТJ,	
		TM,	TN,	TR,	TT,	TZ,	UA,	UG,	US,	UΖ,	VC,	VN,	YU,	ZA,	ZM,	ZW	
	RW:	GH,	GM,	KE,	LS,	MW,	${ m MZ}$,	SD,	SL,	SZ,	${ m TZ}$,	UG,	ZM,	ZW,	ΑM,	AΖ,	
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CA	2506	950			A1		2004	0610	1	CA 2			950		21	0021	125
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	2506						2010										
AU	2002	3652	82		A1		2004	0618	-	AU 2			82		21	0021	125
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EP	1565	416			A1		2005	0824		EP 2			77		21	0021	125
	R:	•	•	•	•	•	ES,	•	•	•	•	•	•	•	•	•	
		•	•	•	•	•	FI,	•	•	•	•	•	•	•	•		
US	2006	0148	976		A1		2006	0706		US 2			06		21	0060.	222
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	7662				В2		2010	0216		^	000	10	200	_		2001	105
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											<						

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT ED Entered STN: 10 Jun 2004

AB A defoaming agent, for cementitious compns., has an excellent defoaming performance and does not segregate when mixed with a high performance air entraining AE water-reducing agent to form a single admixt. or when dild. in H2O at the desired concn., and which has excellent long term storage properties. The defoaming agent is obtained by mixing .gtoreq.1 polyethylene oxide deriv. and a nonionic defoaming agent, where the polyethylene oxide deriv. has at one end a hydrophobic group with a branched structure and/or an unsatd. bond and at the other end an anionic group.

IT 136931-77-6P 478019-97-5P 478019-98-6P 478020-00-7P 478020-01-8P 478020-02-9P

(antifoaming agents for cement compns. contg. nonionic antifoaming agents and polyethylene oxide derivs.)

RN 136931-77-6 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(4-nonylphenoxy)methyl]-2-(2-propen-1-yloxy)ethoxy]-, ammonium salt (1:1) (CA INDEX NAME)

● NH3

RN 478019-97-5 HCAPLUS

CN Poly(oxy-1,2-ethanediy1), .alpha.-sulfo-.omega.-[1-[[(3-phenyl-2-propenyl)oxy]methy1]-2-(2-propenyloxy)ethoxy]-, ammonium salt (9CI) (CA INDEX NAME)

● инз

RN 478019-98-6 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(1-naphthalenyloxy)methyl]-2-(2-propenyloxy)ethoxy]-, ammonium salt (9CI) (CA INDEX NAME)

● NH3

RN 478020-00-7 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-phenyl-2-(2-propenyloxy)ethoxy]-, ammonium salt (9CI) (CA INDEX NAME)

NH3

RN 478020-01-8 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1,2-diphenyl-2-(2-propenyloxy)ethoxy]-, disodium salt (9CI) (CA INDEX NAME)

●2 Na

RN 478020-02-9 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2-[(2-propenyl)phenoxy]ethoxy]-, ammonium salt (9CI) (CA INDEX NAME)

$$CH_2$$
 CH_2
 CH_2

■ NH3

CC 58-1 (Cement, Concrete, and Related Building Materials)
 Section cross-reference(s): 38

31691-97-1P 59764-60-2P 136931-77-6P 171407-73-1P ΙT 478019-97-5P 478019-98-6P 478019-99-7P 478020-01-8P 478020-00-7P 478020-02-9P 478020-04-1P

> (antifoaming agents for cement compns. contg. nonionic antifoaming agents and polyethylene oxide derivs.)

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

APPLICATION NO

RE FORMAT

L39 ANSWER 4 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN 2003:472494 HCAPLUS Full-text ACCESSION NUMBER:

DOCUMENT NUMBER: 139:53794

Reactive liquid polymer crosslinking agent and TITLE:

preparation

INVENTOR(S): Lazar, Warren G.; Clark, James A.

PATENT ASSIGNEE(S): LCB Worldwide Inc., USA SOURCE: PCT Int. Appl., 27 pp.

KIND

CODEN: PIXXD2

DATE

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

PATENT NO

	PAIENI NO.					VIII.		DAIL		APPLICATION NO.						DAIE		
	WO 2003050094			A1 20030619			WO 2002-US38058						20021126					
		W:	ΑE,	AG,	AL,	AM,	ΑT,	AU,	AZ,	BA,	BB,	BG,	BR,	BY,	BZ,	CA,	CH,	
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			LC,	LK,	LR,	LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	
			NO,	NΖ,	OM,	PH,	PL,	PT,	RO,	RU,	SD,	SE,	SG,	SI,	SK,	SL,	ТJ,	
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			EE,	ES,	FI,	FR,	GB,	GR,	ΙE,	ΙΤ,	LU,	MC,	NL,	PT,	SE,	SK,	TR,	
			BF,	ΒJ,	CF,	CG,	CI,	CM,	GΑ,	GN,	GQ,	GW,	ML,	MR,	ΝE,	SN,	TD,	ΤG
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												<						

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

Entered STN: 20 Jun 2003

AΒ A reactive liq. crosslinking agent for use in the prepn. of polymeric substances. The crosslinking agent comprises a substituted (e.g. butylated) 1,3,5-triazine reacted with H2O, an acid alkyl sulfonate and/or phosphonate and a solidifying modifier contg. an hydroxyl functional group. The reactive liq. polymer crosslinking agent has a solids content 20-99% solids. The reactive liq. crosslinking agents (RLPC's) are useful as modifiers in the prepn. of polymeric compds. which are suitable for 1-component self-crosslinking adhesives, coatings and polymers used in optics, textiles, composites, casting and molding. Systems

contg. 1-30% RLPC provide fast single package thermosetting polymeric compds. with enhanced properties such as chem., heat and abrasion resistance.

- IT 1.5214-89-8D, 2-Acrylamido-2-methylpropanesulfonic acid, reaction products with triazine and polyethylene glycol 1.11083-74-0D, reaction products with triazine and diol (reactive liq. polymer crosslinking agent reaction product of substituted triazine, water, sulfonate or phosphonate, and hydroxy compd.)
- RN 15214-89-8 HCAPLUS
- CN 1-Propanesulfonic acid, 2-methyl-2-[(1-oxo-2-propen-1-yl)amino]- (CA INDEX NAME)

- RN 111083-74-0 HCAPLUS
- CN Titanate(3-), [2,2-bis[(2-propen-1-yloxy)methyl]-1-butanolato-.kappa.0][P,P-dioctyl diphosphato(2-)-.kappa.0'']bis[P,P-dioctyl diphosphato(2-)-.kappa.0''']-, hydrogen (1:3) (CA INDEX NAME)

PAGE 1-A

PAGE 2-A

●3 H+

IPCI C07D0251-54 [ICM,7]; C07D0251-66 [ICS,7]; C09K0003-00 [ICS,7]
IPCR C08F0008-30 [I,A]; C09D0167-00 [I,A]; C09J0167-00 [I,A]

CC 37-6 (Plastics Manufacture and Processing)

IT57-50-1D, Sucrose, alkylglycosides, reaction products with triazine and sulfonylzirconate 98-11-3D, Phenylsulfonic acid, reaction products with triazine and diethylene glycol 107-21-1D, Ethylene glycol, reaction products with triazine and sulfonyltitanate 108-78-1D, 2,4,6-Triamino-1,3,5-triazine, reaction products with phenylphosphoric acid 110-63-4D, Butane-1,4-diol, reaction products with triazine and phosphatotitanate 111-46-6D, Diethylene glycol, reaction products with phenylsulfonic acid and triazine 504-63-2D, 1,3-Propylene glycol, reaction products with triazine and sulfonate 629-11-8D, Hexane-1,6-diol, reaction products with triazine and 1571-33-1D, Phenylphosphonic acid, reaction phosphatotitanate products with triazine 5606-17-7D, reaction products with sulfonate and propylene glycol 5606-19-9D, reaction products with polypropylene glycol and sulfate ester 15214-89-8D, 2-Acrylamido-2-methylpropanesulfonic acid, reaction products with triazine and polyethylene glycol 25322-68-3D, Polyethylene glycol, reaction products with triazine and sulfonate 25322-69-4D, Polypropylene glycol, reaction products with triazine and sulfate 89619-91-0D, reaction products with alkylglycosides and sulfonylzirconate 103406-74-2D, reaction products with triazine and ethylene glycol 109766-35-0D, reaction products with triazine and alkylglycosides 111083-74-00, reaction products with triazine and diol 544651-50-5D, reaction products with sulfonate and polyethylene glycol 544651-51-6D, reaction products with phosphate ester and polyethylene glycol 544651-52-7D, reaction products with phosphotitanate and diol

(reactive liq. polymer crosslinking agent reaction product of substituted triazine, water, sulfonate or phosphonate, and hydroxy compd.)

REFERENCE COUNT:

THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L39 ANSWER 5 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2003:222046 HCAPLUS Full-text

5

DOCUMENT NUMBER: 138:260046

TITLE: Method for controlling scale formation and

deposition in aqueous systems

INVENTOR(S): Buentello, Kristin E.; Kessler, Stephen M.; May,

Roger C.; Kaechelin, Julie A.; Chen, Fu; Kolson,

Natalie A.

PATENT ASSIGNEE(S): Betzdearborn Inc., USA

SOURCE: U.S. Pat. Appl. Publ., 10 pp., Cont.-in-part of

U.S. 6,444,747. CODEN: USXXCO

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 4

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 20030052303	A1	20030320	US 2001-878646	20010611

110	6641754	D.O.	20021104	<	
		B2 B1			20010315
CA	2440435	A1	20021010	< CA 2002-2440435 <	20020301
	2440435	C A1			20020301
WO				<	
	CN, CO, GE, GH, LC, LK, NO, NZ, TM, TN,	CR, CU, GM, HR, LR, LS, OM, PH, TR, TT,	CZ, DE, DK, HU, ID, IL, LT, LU, LV, PL, PT, RO, TZ, UA, UG,		FI, GB, GD, KP, KR, KZ, MW, MX, MZ, SK, SL, TJ,
	CH, CY, SE, TR,	DE, DK, BF, BJ,	ES, FI, FR,	SL, SZ, TZ, UG, ZM, Z GB, GR, IE, IT, LU, M CM, GA, GN, GQ, GW, M	MC, NL, PT,
AU	SN, TD, 2002314719		20021015	AU 2002-314719	20020301
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ΕP	1379475	A1	20040114	EP 2002-741643 <	20020301
ΕP			DK, ES, FR,	GB, GR, IT, LI, LU, N	NL, SE, MC,
HU		A2		MK, CY, AL, TR HU 2003-3535 <	20020301
		A3 A		BR 2002-8127	20020301
CN	1496338	А	20040512	< CN 2002-806631 <	20020301
	1262489	C	20060705		
JP	2004528439	Т	20040916	JP 2002-577740 <	20020301
NΖ	528038	А	20050527	NZ 2002-528038 <	20020301
AT	435839	Т	20090715	AT 2002-741643	20020301
PT	1379475	E	20090901	PT 2002-741643	20020301
ES	2326960	Т3	20091022	ES 2002-741643	20020301
PL	204918	В1	20100226	PL 2002-363968	20020301
TW	300060	В	20080821	< TW 2002-104266 <	20020307
IN	2003KN00910	А	20050708	IN 2003-KN910	20030715
US	20040039144	A1	20040226	< US 2003-646278 <	20030822
	7094852 904789	B2 B1	20060822 20090625	KR 2003-7011900 <	20030909

10/596,747

NO 2003004049	А	20031104	NO	2003-4049		20030912
MX 2003008342	А	20041015	MX	2003-8342		20030912
AU 2007202743	A1	20070705	AU	2007-202743		20070613
AU 2007202743	В2	20081120				
PRIORITY APPLN. INFO.:			US	2001-808679	Α2	20010315
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				<		
			US	2001-878646	Α	20010611
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			AU	2002-314719	A3	20020301
				<		
			WO	2002-US6370	M	20020301
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ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT ED Entered STN: 21 Mar 2003

GΙ

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$$[E]_{e}$$
* * CH_{2} - $CH_{$

AB A method for inhibiting the formation and deposition of scale forming moieties in aq. systems by adding water-sol. or water-dispersible polymers is disclosed. The method comprises adding to an aq. system a polymer having repeat units characterized by I, wherein E is the repeat unit remaining after polymn. of an ethylenically unsatd. compd.; preferably, a carboxylic acid, sulfonic acid, phosphonic acid, or amide form thereof or mixts. thereof, R1 is H or lower (C1-C4) alkyl, G is -CH2-or -CHCH3-; R2 is CH2-CH2-On or CH2-CHCH3-n where n ranges from about 1 to 100, preferably about 1 to 20, X is an anionic radical selected from the group consisting

of SO3, PO3, or COO; Z is H or hydrogens or any water sol. cationic moiety which counterbalances the valence of the anionic radical X, including but not limited to Na, K, Ca, or NH4, F, when present, is a repeat unit having the structure of II, wherein X and Z are the same as in Formula I. R4 is H or lower (C1-C4) alkyl, R5 is hydroxy substituted alkyl or alkylene having from about 1 to 6 of carbon atoms. 330666-77-8P 452311-66-9P ΙT 171439-08-0P 452311-67-0P 452311-68-1P 452311-69-2P oxide-allyloxy-2-hydroxypropane-3-sulfonic acid graft terpolymer, ammonium sulfate 502546-11-49 (method for controlling scale formation and deposition in aq. systems) RN171439-08-0 HCAPLUS Poly(oxy-1,2-ethanediy1), .alpha.-phosphono-.omega.-(2-propen-1-yloxy)-CN (CA INDEX NAME)

$$H_2C$$
 = $CH_2 - CH_2 - CH_2 - CH_2 - O$ n PO_3H_2

RN 330666-77-8 HCAPLUS
CN 2-Propenoic acid, polymer with
.alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl)
ammonium salt (1:1) (CA INDEX NAME)

CM 1

CRN 55866-85-8
CMF (C2 H4 O)n C3 H6 O4 S . H3 N
CCI PMS

$$H_2C$$
 $=$ CH_2-CH_2-O $=$ CH_2-CH_2-O $=$ SO_3H $=$ NH_3

CM 2

CRN 79-10-7

CMF C3 H4 O2

452311-66-9 HCAPLUS RN 2-Propenoic acid, polymer with 2-hydroxy-3-(2-propen-1-yloxy)-1-propanesulfonic acid and .alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl) ammonium salt (1:1) (CA INDEX NAME) CM1 CRN 94928-31-1 CMF C6 H12 O5 S ОН HO3S-CH2-CH-CH2-O-CH2-CH=CH2 CM CRN 55866-85-8 CMF (C2 H4 O)n C3 H6 O4 S . H3 N CCI PMS H₂C=CH-CH₂-O-CH₂-CH₂-O-I_n SO₃H ● NH3 CM 3 CRN 79-10-7 CMF C3 H4 O2 HO-C-CH-CH2 452311-67-0 HCAPLUS RN 2-Propenoic acid, 2-methyl-, polymer with 2-propenoic acid and .alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl)

ammonium salt (1:1) (CA INDEX NAME)

СМ

CM 2

CRN 79-41-4 CMF C4 H6 O2

CM 3

CRN 79-10-7 CMF C3 H4 O2

RN 452311-68-1 HCAPLUS
CN 2-Propenoic acid, polymer with
2-methyl-2-[(1-oxo-2-propen-1-yl)amino]-1-propanesulfonic acid and
.alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl)
ammonium salt (1:1) (CA INDEX NAME)

CM 1

CRN 55866-85-8

CMF (C2 H4 O)n C3 H6 O4 S . H3 N

CCI PMS

CM 2

CRN 15214-89-8 CMF C7 H13 N O4 S

CM 3

CRN 79-10-7 CMF C3 H4 O2

CM 1

CRN 171439-08-0

CMF (C2 H4 O)n C3 H7 O4 P

CCI PMS

$$H_2C$$
 = $CH_2 - CH_2 - CH_2 - CH_2 - CH_2 - O$ n $PO3H_2$

CMF H2 O4 S

CM 2

CRN 502546-08-9

CMF (C6 H12 O5 S . C3 H4 O2 . C2 H4 O) x

CCI PMS

CM 3

CRN 94928-31-1

CMF C6 H12 O5 S

CM 4

CRN 79-10-7

CMF C3 H4 O2

CM 5

CRN 75-21-8

CMF C2 H4 O



RN 502546-11-4 HCAPLUS

CN 2-Propenoic acid, polymer with

2-methyl-2-[(1-oxo-2-propenyl)amino]-1-propanesulfonic acid and oxirane, hydrogen sulfate (ester), graft, ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 7664-93-9 CMF H2 O4 S

CM 2

CRN 164578-98-7

CMF (C7 H13 N O4 S . C3 H4 O2 . C2 H4 O) \times

CCI PMS

CM 3

CRN 15214-89-8 CMF C7 H13 N O4 S

CM 4

CRN 79-10-7 CMF C3 H4 O2

CM 5

CRN 75-21-8 CMF C2 H4 O

ED Entered STN: 13 Dec 2002



INCL 252175000 IPCI C02F0005-02 [ICM,7] IPCR F25D0017-02 [I,A]; C02F0005-00 [I,A]; C02F0005-10 [I,A]; C02F0005-12 [I,A]; C02F0005-14 [I,A]; C08F0216-14 [I,A]; C23F0011-16 [I,A]; C23F0011-167 [I,A] NCL 252/175.000; 252/180.000; 252/181.000; 524/807.000; 526/287.000 CC 61-8 (Water) ΤТ 171439-08-0P 330666-77-8P 452311-66-9P 452311-67-0P 452311-68-1P 452311-69-2P **452311-70-5P** 502546-07-8P, Acrylic acid-ethylene oxide graft copolymer, ammonium sulfate 502546-09-02, Acrylic acid-ethylene oxide-allyloxy-2-hydroxypropane-3-sulfonic acid graft terpolymer, ammonium sulfate 502546-11-49 502546-12-5P 502546-13-6P (method for controlling scale formation and deposition in aq. systems) OS.CITING REF COUNT: 4 THERE ARE 4 CAPLUS RECORDS THAT CITE THIS RECORD (4 CITINGS) L39 ANSWER 6 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2002:945818 HCAPLUS Full-text 138:28220 DOCUMENT NUMBER: TITLE: Antifoaming agents for cement compositions showing segregation resistance when being mixed with high-performance air entraining water reducing agents or water INVENTOR(S): Lorentz, Claus; Yaguchi, Minoru; Sugiyama, Tomomi; Albrecht, Gerhard NMB K. K., Japan PATENT ASSIGNEE(S): Jpn. Kokai Tokkyo Koho, 9 pp. SOURCE: CODEN: JKXXAF DOCUMENT TYPE: Patent LANGUAGE: Japanese FAMILY ACC. NUM. COUNT: 1 PATENT INFORMATION: APPLICATION NO. PATENT NO. KIND DATE ____ 20010531 A 20021213 JP 2001-164388 JP 2002356356 <--PRIORITY APPLN. INFO.: JP 2001-164388 20010531

<--

10/596,747

AB The antifoaming agents contain nonionic antifoaming agents and polyethylene oxide derivs. having one anionic terminals and branched and/or unsatd. hydrophobic terminals on the other end.

IT 136931-77-6P 478019-97-5P 478019-98-6P 478020-00-7P 478020-01-8P 478020-02-9P

(antifoaming agents for cement compns. contg. nonionic antifoaming agents and polyethylene oxide derivs.)

RN 136931-77-6 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(4-nonylphenoxy)methyl]-2-(2-propen-1-yloxy)ethoxy]-, ammonium salt (1:1) (CA INDEX NAME)

● NH3

RN 478019-97-5 HCAPLUS

CN Poly(oxy-1,2-ethanediy1), .alpha.-sulfo-.omega.-[1-[[(3-phenyl-2-propenyl)oxy]methy1]-2-(2-propenyloxy)ethoxy]-, ammonium salt (9CI) (CA INDEX NAME)

● инз

RN 478019-98-6 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(1-naphthalenyloxy)methyl]-2-(2-propenyloxy)ethoxy]-, ammonium salt (9CI) (CA INDEX NAME)

● NH3

RN 478020-00-7 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-phenyl-2-(2-propenyloxy)ethoxy]-, ammonium salt (9CI) (CA INDEX NAME)

● NH3

RN 478020-01-8 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1,2-diphenyl-2-(2-propenyloxy)ethoxy]-, disodium salt (9CI) (CA INDEX NAME)

$$\label{eq:h2C} \begin{array}{c} \text{Ph} & \text{Ph} \\ \text{H2C} = \text{CH} - \text{CH}_2 - \text{O} - \text{CH} - \text{CH} - \text{O} - \text{CH}_2 - \text{CH}_2 - \text{O} - \text{O} \\ & & & & & & & & & & & & \\ \end{array}$$

•2 Na

RN 478020-02-9 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2-[(2-propenyl)phenoxy]ethoxy]-, ammonium salt (9CI) (CA INDEX NAME)

$$CH_2$$
 CH_2
 CH_2

● NH3

CC 58-1 (Cement, Concrete, and Related Building Materials)

Section cross-reference(s): 38

IT 31691-97-1P 59764-60-2P 136931-77-6P 171407-73-1P

478019-97-5P 478019-98-6P 478019-99-7P **478020-00-7P 478020-01-8P 478020-02-9P**

478020-04-1P

(antifoaming agents for cement compns. contg. nonionic antifoaming agents and polyethylene oxide derivs.)

L39 ANSWER 7 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2002:777854 HCAPLUS Full-text

DOCUMENT NUMBER: 137:299503

TITLE: Method for controlling scale formation and

deposition in aqueous systems

INVENTOR(S): Chen, Fu; Kolson, Natalie A.; Buentello, Kristin

E.; Kaechelin, Julie A.; Kessler, Stephen M.; May,

Roger C.

PATENT ASSIGNEE(S): Betzdearborn Inc., USA SOURCE: PCT Int. Appl., 39 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 4

PATENT INFORMATION:

PATENT NO	TENT NO. KIND DATE APPLICATION NO.							DATE							
WO 2002079106 A1			A1 20021010			WO 2002-US6370 <						20020301			
RW: (AE, A CN, GE, G LC, I NO, I TM, G GH, G CH, G SE, SN, S	CO, GH, LK, NZ, TN, GM, CY,	CR, GM, LR, OM, TR, KE, DE, BF,	CU, HR, LS, PH, TT, LS, DK,	CZ, HU, LT, PL, TZ, MW, ES,	DE, ID, LU, PT, UA, MZ, FI,	DK, IL, LV, RO, UG, SD, FR,	DM, IN, MA, RU, UZ, SL, GB,	DZ, IS, MD, SD, VN, SZ, GR,	BG, EC, JP, MG, SE, YU, TZ, IE,	BR, EE, KE, MK, SG, ZA, UG, IT,	ES, KG, MN, SI, ZM, ZM, LU,	FI, KP, MW, SK, ZW ZW, MC,	GB, KR, MX, SL, AT, NL,	GD, KZ, MZ, TJ, BE, PT,

10/596,747

US	6444747	В1	20020903	US 2001-808679		20010315
US	20030052303	A1	20030320	< US 2001-878646 <		20010611
US	6641754	В2	20031104			
CA	2440435	A1	20021010	CA 2002-2440435		20020301
CZ	2440435	С	20100608	<		
	2002314719	A1	20021015	AU 2002-314719		20020301
	000001.151.0	- 0	000=0.00	<		
	2002314719 1379475	B2 A1	20070426 20040114	EP 2002-741643		20020301
111	13/34/3	711	20040114	<		20020301
EP	1379475	B1				
				GB, GR, IT, LI, LU, MK, CY, AL, TR	NL, SE	E, MC,
HU	2003003535	A2		HU 2003-3535		20020301
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	2003003535 2002008127	A3 A	20080328 20040323			20020301
DIX	2002000127	A	20040323	SR 2002 0127		20020301
JP	2004528439	Т	20040916	JP 2002-577740 <		20020301
NΖ	528038	А	20050527	NZ 2002-528038		20020301
АТ	435839	Т	20090715	< AT 2002-741643		20020301
				<		
IN	2003KN00910	А	20050708	IN 2003-KN910 <		20030715
KR	904789	В1	20090625	KR 2003-7011900		20030909
				<		
NO	2003004049	А	20031104	NO 2003-4049		20030912
MX	2003008342	А	20041015	< MX 2003-8342		20030912
				<		
AU	2007202743	A1	20070705	AU 2007-202743		20070613
AU	2007202743	В2	20081120	<		
PRIORITY	Y APPLN. INFO.:			US 2001-806679	А	20010315
				<	-70	00010015
				US 2001-808679 <	A	20010315
				US 2001-878646	А	20010611
				<		0000000
				AU 2002-314719 <	АЗ	20020301
				WO 2002-US6370	W	20020301
				<		

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT Entered STN: 11 Oct 2002

AΒ

A water sol. or water dispersible polymer compn., useful as scale and/or corrosion inhibitors in aq. systems has the formula [E]c[CH2CR1(GOR2XZ)]d[F]e wherein E is the repeating unit remaining after polymn. of an ethylenically unsatd. compd., preferably, a carboxylic acid, sulfonic acid, phosphonic acid, or amide form thereof or mixts. thereof; R1 is H or lower C1-4 alkyl; G is CH2 or CHMe; R2 is (CH2CH2O)n or (CH2CHMeO)m where n and m range from about 1 to 100, preferably n

10/596.747

is greater than 10 and m ranges from about 1 to 20; X is an anionic radical selected from the group consisting of SO3, PO3, or CO2; Z is H or hydrogens or any water sol. cationic moiety which counterbalances the valence of the anionic radical X, including but not limited to Na, K, Ca, or NH4; F, when present, is a repeating unit having the structure of formula [CH2CR4(CH2OR5XZ)] wherein X and Z are the same as above, R4 is H or a lower C1-4 alkyl, and R5 is a hydroxy-substituted alkyl or alkylene having from about 1 to 6 carbon atoms. This water sol. or water dispersible polymer compn. is useful: as scale deposit control and corrosion inhibition agents in water treatment or gas scrubbing processes, in pulp and paper manufg. processes, in pretreating of metals; as rheol. modifiers for concrete and cement additives; as cleaning agents for membranes; and as hydrophilic modifier components in personal care, cosmetic and pharmaceutical formulations.

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ΙT
     171439-08-0P, Poly(oxy-1,2-ethanediyl),
     .alpha.-phosphono-.omega.-(2-propenyloxy)- 330666-77-8P,
     2-Propenoic acid, polymer with
     .alpha.-sulfo-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl) ammonium
            452311-66-9P, 2-Propenoic acid, polymer with
     2-hydroxy-3-(2-propenyloxy)-1-propanesulfonic acid and
     .alpha.-sulfo-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl) ammonium
           452311-67-0P, 2-Propenoic acid, 2-methyl-, polymer
     with 2-propenoic acid and .alpha.-sulfo-.omega.-(2-
     propenyloxy) poly (oxy-1, 2-ethanediyl) ammonium salt
     452311-68-1P, 2-Propenoic acid, polymer with
     2-methyl-2-[(1-oxo-2-propenyl)amino]-1-propanesulfonic acid and
     .alpha.-sulfo-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl) ammonium
           452311-69-29, 2-Propenoic acid, polymer with
     .alpha.-phosphono-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl)
     452311-70-5P, 2-Propenoic acid, polymer with
     .alpha.-sulfo-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl)
     452311-71-6P
        (water sol. or water dispersible polymers as scale and/or corrosion
        inhibitors in aq. systems)
RN
     171439-08-0 HCAPLUS
     Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-(2-propen-1-yloxy)-
CN
       (CA INDEX NAME)
```

$$\text{H}_2\text{C} \textcolor{red}{==} \text{CH} \textcolor{blue}{-} \text{CH}_2 \textcolor{blue}{-} \text{O} \textcolor{blue}{-} \textcolor{blue}{-} \text{CH}_2 \textcolor{blue}{-} \text{O} \textcolor{blue}{-} \textcolor{blue}{-} \textcolor{blue}{-} \text{PO}_3\text{H}_2$$

CM 2

CRN 79-10-7 CMF C3 H4 O2

RN 452311-66-9 HCAPLUS

CN 2-Propenoic acid, polymer with 2-hydroxy-3-(2-propen-1-yloxy)-1-propanesulfonic acid and .alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl) ammonium salt (1:1) (CA INDEX NAME)

CM 1

CRN 94928-31-1 CMF C6 H12 O5 S

CM 2

CRN 55866-85-8 CMF (C2 H4 O)n C3 H6 O4 S . H3 N CCI PMS

● NH3

RN 452311-67-0 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, polymer with 2-propenoic acid and .alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl) ammonium salt (1:1) (CA INDEX NAME)

CM 1

CRN 55866-85-8

CMF (C2 H4 O)n C3 H6 O4 S . H3 N $\,$

CCI PMS

$$H_2C$$
 = $CH_2 - CH_2 - CH_2 - CH_2 - CH_2 - O$ n SO_3H

● NH3

CM 2

CRN 79-41-4 CMF C4 H6 O2

CM 3

CRN 79-10-7 CMF C3 H4 O2

452311-68-1 HCAPLUS RNCN 2-Propenoic acid, polymer with

2-methyl-2-[(1-oxo-2-propen-1-yl)amino]-1-propanesulfonic acid and .alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl)

ammonium salt (1:1) (CA INDEX NAME)

CM1

CRN 55866-85-8

CMF (C2 H4 O)n C3 H6 O4 S . H3 N

CCI PMS

CM

CRN 15214-89-8 CMF C7 H13 N O4 S

3 CM

CRN 79-10-7 CMF C3 H4 O2

```
RN
     452311-69-2 HCAPLUS
CN
     2-Propenoic acid, polymer with
     .alpha.-phosphono-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl)
     (9CI) (CA INDEX NAME)
     CM
          1
     CRN 171439-08-0
     CMF
          (C2 H4 O)n C3 H7 O4 P
     CCI PMS
H_2C = CH - CH_2 - O = CH_2 - CH_2 - O = DO3H_2
     CM
     CRN 79-10-7
     CMF C3 H4 O2
     452311-70-5 HCAPLUS
RN
     2-Propenoic acid, polymer with
CN
     .alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl)
                                                                            (CA
     INDEX NAME)
          1
     CM
     CRN 201605-73-4
     CMF (C2 H4 O)n C3 H6 O4 S
     CCI PMS
H<sub>2</sub>C=CH-CH<sub>2</sub>-O-CH<sub>2</sub>-CH<sub>2</sub>-O-In SO<sub>3</sub>H
          2
     CM
     CRN 79-10-7
```

CMF C3 H4 O2

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IPCI C02F0005-10 [ICM,7]; C08F0216-14 [ICS,7]
IPCR F25D0017-02 [I,A]; C02F0005-00 [I,A]; C02F0005-10 [I,A]; C02F0005-12
     [I,A]; C02F0005-14 [I,A]; C08F0216-14 [I,A]; C23F0011-16 [I,A];
     C23F0011-167 [I,A]
     61-8 (Water)
CC
     Section cross-reference(s): 35, 43, 58, 59, 62, 63
     171439-08-0P, Poly(oxy-1,2-ethanediyl),
IΤ
     .alpha.-phosphono-.omega.-(2-propenyloxy)- 330666-77-8P,
     2-Propenoic acid, polymer with
     .alpha.-sulfo-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl) ammonium
          452311-66-9P, 2-Propenoic acid, polymer with
     2-hydroxy-3-(2-propenyloxy)-1-propanesulfonic acid and
     .alpha.-sulfo-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl) ammonium
     salt 452311-67-0P, 2-Propenoic acid, 2-methyl-, polymer
```

10/596,747

with 2-propenoic acid and .alpha.-sulfo-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl) ammonium salt 452311-68-19, 2-Propenoic acid, polymer with 2-methyl-2-[(1-oxo-2-propenyl)amino]-1-propanesulfonic acid and .alpha.-sulfo-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl) ammonium salt 452311-69-29, 2-Propenoic acid, polymer with .alpha.-phosphono-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl) 452311-70-59, 2-Propenoic acid, polymer with .alpha.-sulfo-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl) 452311-71-69

(water sol. or water dispersible polymers as scale and/or corrosion inhibitors in aq. systems)

OS.CITING REF COUNT: 1 THERE ARE 1 CAPLUS RECORDS THAT CITE THIS

RECORD (1 CITINGS)

REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L39 ANSWER 8 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2002:669688 HCAPLUS Full-text

DOCUMENT NUMBER: 137:202027

TITLE: Water soluble polymers
INVENTOR(S): Chen, Fu; Kolson, Natalie A.

PATENT ASSIGNEE(S): Betzdearborn Inc., USA

SOURCE: U.S., 7 pp.
CODEN: USXXAM

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 4

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 6444747	B1	20020903	US 2001-808679	20010315
US 20030052303	A1	20030320	US 2001-878646	20010611
US 6641754 CA 2440435		20031104 20021010	< CA 2002-2440435	20020301
CA 2440435 WO 2002079106	•	20100608 20021010	< WO 2002-US6370	20020301
CN, CO, GE, GH, LC, LK, NO, NZ, TM, TN,	CR, CU, CZ, GM, HR, HU, LR, LS, LT, OM, PH, PL, TR, TT, TZ,	DE, DK, I ID, IL, I LU, LV, N PT, RO, I UA, UG, U	<pre></pre>	FI, GB, GD, KP, KR, KZ, AW, MX, MZ, SK, SL, TJ,
CH, CY,	DE, DK, ES, BF, BJ, CF,	FI, FR, (GB, GR, IE, IT, LU, N CM, GA, GN, GQ, GW, N	MC, NL, PT,
- , ,		20021015	AU 2002-314719	20020301
AU 2002314719	В2	20070426	`	

10/596,747

EP	1379475	5		A1		20040	0114	Ι	EΡ	200			43			20	0203	301
EP		BE,			DK,		FR,					LI,	LU,	NL,	SE	,	MC,	
HU	PT 2003003	E, IE, 8535	SI,	LT, A2		FI, 2004						3535				20	0203	301
	2003003 1496338			A3 A		20080			CN	200	2-8	3066	31			20	0200	301
	1262489 2004528			C T		20060		ı	JP	200		5777	40			20	0203	301
NZ	528038			А		20050	0527	1	ΝZ	200	2-5	 5280	38			20	0203	301
AT	435839			Т		20090	0715	Ī	ΑT	200	2-7	 7416	43			20	0203	301
PT	1379475	5		E		20090	0901	1	PΤ	200	2-7	 7416	43			20	0203	301
ES	2326960)		Т3		2009	1022	Ι	ES			7416	43			20	0203	301
PL	204918			В1		2010	0226]	ΡL			3639	68			20	0203	301
TW	300060			В		20080	0821	ŗ	ΓW	200		042	66			20	0203	307
IN	2003KN0	0910		А		20050	0708	-	IN	200		KN91	0			20	030	715
US	2004003	39144		A1		20040	0226	Ţ	US	200	<- 3-6 <-	5462	78			20	0308	822
	7094852 904789)		B2 B1		20060 20090		I	KR	200	3-7	7011	900			20	0309	909
NO	2003004	1049		А		2003	1104	I	NO	200	3-4	1049				20	0309	912
MX	2003008	342		А		2004	1015	1	MX	200	3-8	342				20	0309	912
AU	2007202	2743		A1		20070	0705	Ž	AU	200		2027	43			20	070	613
AU PRIORITY	2007202 APPLN.		. •	В2		2008	1120	ī	US	200		3066	79		Α	20	0103	315
11(101(111		11110	• •								<-						0103	
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											<-						0203	
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								`	W	200	<-		, 0		4.4	۷ ر	,020.) () <u>T</u>

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT ED Entered STN: 05 Sep 2002 GI

* — [E]c — * *
$$\begin{array}{c} & & \\ &$$

AB A water sol. or water dispersible polymer compn., useful as deposit control and corrosion inhibition agents in water treatment, pulp and paper manufg. processes, and in pretreating of metals; as rheol. modifiers for concrete and cement additives; as cleaning agents for membranes; and as hydrophilic modifier components in personal care, cosmetic and pharmaceutical formulations, has the repeat units of I: Wherein E is the repeat unit remaining after polymn. of an ethylenically unsatd. compd.; preferably, a carboxylic acid, sulfonic acid, phosphonic acid, or amide thereof or mixts. thereof; R1 is H or C1-C4 alkyl; G is -CH2- or -CHCH3-; R2 is polyoxyethylene or polyoxypropylene where n is 1-100, preferably 1-20; X is an anionic radical selected from the group consisting of SO3, PO3, or CO2; Z is H or hydrogens or any water sol. cationic moiety which counterbalances the valence of the anionic radical X, including but not limited to Na, K, Ca, or NH4; F, when present, is a repeat unit having the structure of II: wherein X and Z are the same as in I; R4 is H or C1-C4 alkyl; R5 is hydroxy substituted alkyl or C1-C6 alkylene.

IT 330666-77-8p 452311-66-9p 452311-67-0p 452311-68-1p 452311-69-2p 452311-70-5p

(water sol. polymers)

RN 330666-77-8 HCAPLUS

2-Propenoic acid, polymer with .alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl) ammonium salt (1:1) (CA INDEX NAME)

CM 1

CN

CRN 55866-85-8 CMF (C2 H4 O)n C3 H6 O4 S . H3 N CCI PMS

● NH3

CRN 79-10-7 CMF C3 H4 O2

RN 452311-66-9 HCAPLUS

CN 2-Propenoic acid, polymer with 2-hydroxy-3-(2-propen-1-yloxy)-1-propanesulfonic acid and .alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl) ammonium salt (1:1) (CA INDEX NAME)

CM 1

CRN 94928-31-1 CMF C6 H12 O5 S

CM 2

CRN 55866-85-8 CMF (C2 H4 O)n C3 H6 O4 S . H3 N CCI PMS

● NH3

CM 3

CRN 79-10-7 CMF C3 H4 O2

RN 452311-67-0 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, polymer with 2-propenoic acid and .alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl) ammonium salt (1:1) (CA INDEX NAME)

CM 1

CRN 55866-85-8

CMF (C2 H4 O)n C3 H6 O4 S . H3 N

CCI PMS

$$H_2C$$
 $=$ CH_2-CH_2-O $=$ CH_2-CH_2-O $=$ SO_3H $=$ NH_3

CM 2

CRN 79-41-4 CMF C4 H6 O2

CM 3

CRN 79-10-7 CMF C3 H4 O2

RN 452311-68-1 HCAPLUS

CN 2-Propenoic acid, polymer with 2-methyl-2-[(1-oxo-2-propen-1-yl)amino]-1-propanesulfonic acid and .alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl) ammonium salt (1:1) (CA INDEX NAME)

CM 1

CRN 55866-85-8 CMF (C2 H4 O)n C3 H6 O4 S . H3 N CCI PMS

● NH3

CM 2

CRN 15214-89-8 CMF C7 H13 N O4 S

CM 3

CRN 79-10-7 CMF C3 H4 O2

CRN 171439-08-0 CMF (C2 H4 O)n C3 H7 O4 P

CCI PMS

$$H_2C$$
 = $CH_2 - CH_2 - CH_2 - CH_2 - O$ n $PO3H_2$

CM 2

CRN 79-10-7 CMF C3 H4 O2

RN 452311-70-5 HCAPLUS

CN 2-Propenoic acid, polymer with

.alpha.-sulfo-.omega.-(2-propen-1-yloxy)poly(oxy-1,2-ethanediyl) (CA INDEX NAME)

CM 1

CRN 201605-73-4

CMF (C2 H4 O)n C3 H6 O4 S

CCI PMS

CM 2

CRN 79-10-7

CMF C3 H4 O2

```
452311-71-6
ΙT
         (water sol. polymers)
RN
     452311-71-6 HCAPLUS
     2-Propenoic acid, 2-methyl-, polymer with
CN
     .alpha.-sulfo-.omega.-(2-propenyloxy)poly(oxy-1,2-ethanediyl) ammonium
     salt (9CI) (CA INDEX NAME)
     CM
           1
     CRN 55866-85-8
           (C2 H4 O)n C3 H6 O4 S . H3 N
     CMF
     CCI
H<sub>2</sub>C=CH-CH<sub>2</sub>-O-CH<sub>2</sub>-CH<sub>2</sub>-O-In SO<sub>3</sub>H
                     ● NH3
     CM
     CRN 79-41-4
     CMF C4 H6 O2
CH2
||
Me-C-CO2H
     171439-08-0P
ΙT
         (water sol. polymers)
     171439-08-0 HCAPLUS
RN
CN
     Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-(2-propen-1-yloxy)-
        (CA INDEX NAME)
H_2C \longrightarrow CH_2-CH_2-O \longrightarrow CH_2-CH_2-O \longrightarrow DO3H_2
INCL 524807000
IPCI C08L0041-00 [ICM,7]; C08L0043-00 [ICS,7]; C08F0220-04 [ICS,7];
     C08F0220-64 [ICS,7]; C08F0228-02 [ICS,7]
IPCR C02F0005-10 [I,A]; C02F0005-12 [I,A]; C02F0005-14 [I,A]; C08F0216-14
     [I,A]
NCL 524/807.000; 524/817.000; 524/832.000; 526/287.000; 526/318.410;
```

10/596,747

526/320.000

CC 37-3 (Plastics Manufacture and Processing)

Section cross-reference(s): 43, 58, 62, 63

IT 330666-77-8P 452311-66-9P 452311-67-0P

452311-68-1P 452311-69-2P 452311-70-5P

(water sol. polymers)

IT 452311-71-6

(water sol. polymers)

IT 171439-08-0P

(water sol. polymers)

OS.CITING REF COUNT: 4 THERE ARE 4 CAPLUS RECORDS THAT CITE THIS

RECORD (4 CITINGS)

REFERENCE COUNT: 21 THERE ARE 21 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L39 ANSWER 9 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2001:603583 HCAPLUS Full-text

DOCUMENT NUMBER: 135:167504

TITLE: Reactive surfactant compositions and manufacture

of polymer emulsions using them

INVENTOR(S): Ishikawa, Yoshinobu; Sawada, Hiroki; Ishii, Yasuo

PATENT ASSIGNEE(S): Kao Corp., Japan

SOURCE: Jpn. Kokai Tokkyo Koho, 9 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PA:	TENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP	2001226439	A	20010821	JP 2000-33804	20000210
-	3420733 1129770	B2 A1	20030630 20010905	EP 2001-102796	20010209
EP	1129770 R: AT, BE, CH, PT, IE, SI,			B, GR, IT, LI, LU, NI	, SE, MC,
US		A1		US 2001-779314 <	20010209
ES	2230186	Т3	20050501	ES 2001-102796 <	20010209
PRIORITY	Y APPLN. INFO.:			JP 2000-33804 <	A 20000210

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

ED Entered STN: 21 Aug 2001

AB The compns. essentially contain reactive surfactants (a) [crit. micelle concn. (CMC) 0.007-0.4 mol/L] having one reactive group selected from CHR1:CR2(Y)p (R1, R2 = H, CH2X; X = H, substituent; Y = CO, CH2; p = 0, 1) and one anionic hydrophilic group and reactive surfactants (b) (CMC 1 .times. 10-5-0.007 mol/L) having one reactive group above mentioned and one anionic hydrophilic group [a/b = 5/95 to 60/40 (by mol)]. Thus, a polymer emulsion prepd. from 71.5 g H2O, 1g of a compn. contg. a 12.9:87.1 mixt. of CH2:C(CO2K)CH2CO2-n-C6H13 and

10/596,747

CH2:C(CO2K)CH2CO2-n-C12H25, 50g styrene, and 7.5g 2% aq. soln. of K2S2O8 showed good polymn. stability and av. particle size 85.1 nm.

IT 354552-66-2P 354552-68-4P 354583-13-4P

354583-14-5P 354583-15-6P 354583-16-7P

354583-17-8P 354583-18-9P

(reactive surfactant compns. with controlled CMC for polymer emulsions)

RN 354552-66-2 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with dodecyl 2-hydroxy-3-(2-propenyloxy)propyl hydrogen phosphate monosodium salt and .alpha.-sulfo-.omega.-[1-[[4-(1,1-dimethylethyl)phenoxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 354552-65-1 CMF C18 H37 O6 P . Na

Na

CM 2

CRN 354552-64-0

CMF (C2 H4 O)n C16 H24 O6 S . H3 N

CCI PMS

■ NH3

CM 3

CRN 141-32-2

CMF C7 H12 O2

RN 354552-68-4 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with dodecyl 2-hydroxy-3-(2-propenyloxy)propyl hydrogen phosphate monosodium salt and 2-hydroxy-3-(2-propenyloxy)propyl octyl hydrogen phosphate monosodium salt (9CI) (CA INDEX NAME)

CM 1

CRN 354552-67-3

CMF C14 H29 O6 P . Na

● Na

CM 2

CRN 354552-65-1

CMF C18 H37 O6 P . Na

Na

CM 3

CRN 141-32-2

CMF C7 H12 O2

RN 354583-13-4 HCAPLUS

CN Butanedioic acid, sulfo-, C-[2-hydroxy-3-(2-propenyloxy)propyl]
C-octyl ester, monosodium salt, polymer with butyl 2-propenoate,
methyl 2-methyl-2-propenoate, 2-propenoic acid and
.alpha.-sulfo-.omega.-[1-[(4-isononylphenoxy)methyl]-2-(2propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA
INDEX NAME)

CM 1

CRN 354583-12-3 CMF (C2 H4 O)n C21 H34 O6 S . H3 N CCI IDS, PMS

$$\begin{array}{c|c} O & \begin{array}{c} CH_2-CH_2-O & \\ \hline \end{array} & SO_3H \\ \hline \\ (iso-C9H_19) & \end{array}$$

● NH3

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 80-62-6 CMF C5 H8 O2

RN 354583-14-5 HCAPLUS

CN Butanedioic acid, sulfo-, C-dodecyl

C-[2-hydroxy-3-(2-propenyloxy)propyl] ester, monosodium salt, polymer with butyl 2-propenoate, methyl 2-methyl-2-propenoate, 2-propenoic acid and .alpha.-sulfo-.omega.-[1-[[4-(1,1-dimethylethyl)phenoxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 354552-64-0

CMF (C2 H4 O)n C16 H24 O6 S . H3 N

CCI PMS

$$\begin{array}{c|c} \text{HO}_{3}\text{S} & \hline & \text{O}_{-\text{CH}_2-\text{CH}_2} \\ \text{H}_{2}\text{C} & \hline & \text{CH}_{-\text{CH}_2-\text{O}_{-\text{CH}_2}-\text{CH}_{-\text{CH}_2}-\text{O}_{-\text{CH}_2}-\text{CH}_{-\text{CH}_2-\text{O}_{-\text{CH}_2}-\text{CH}_{-\text{CH}_2}-\text{CH}_{-\text{CH}_2}-\text{CH}_{-\text{CH}_2-\text{CH}_2-\text{CH}_2-\text{CH}_2-\text{CH}_{-\text{CH}_2}-\text{CH}_{-\text{CH}_2}-\text{CH}_{-\text{CH}_2$$

● NH3

CM 2

CRN 141-32-2

CMF C7 H12 O2

CM 3

CRN 80-62-6

CMF C5 H8 O2

CCI IDS

$$HO-(CH_2)_{11}-Me$$

CN 2-Propenoic acid, 2-methyl-, methyl ester, polymer with butyl 2-propenoate, 2-propenoic acid, alpha.-sulfo-.omega.-[1-[[4-(1,1-dimethylethyl)phenoxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt and alpha.-sulfo-.omega.-[4-isononyl(1-propenyl)phenoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 354583-10-1 CMF (C2 H4 O)n C18 H28 O4 S . H3 N CCI IDS, PMS

$$\begin{array}{c|c} O & \hline \\ CH_2 - CH_2 - O \\ \hline \\ n \end{array} SO_3H$$

D1-CH — CH-Me

● NH3

CM 2

CRN 354552-64-0 CMF (C2 H4 O)n C16 H24 O6 S . H3 N CCI PMS

$$\begin{array}{c|c} \text{HO}_{3}\text{S} & \hline & \text{O-CH}_2\text{-CH}_2 \\ \hline & \text{H}_2\text{C} = \text{CH-CH}_2\text{-O-CH}_2\text{-CH-CH}_2\text{-O} \\ \end{array}$$

● NH3

CM 3

CRN 141-32-2 CMF C7 H12 O2

CM 4

CRN 80-62-6 CMF C5 H8 O2

CM 5

CRN 79-10-7 CMF C3 H4 O2

RN 354583-16-7 HCAPLUS

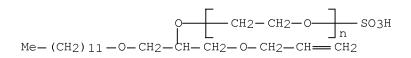
CN Butanedioic acid, sulfo-, C-[2-hydroxy-3-(2-propenyloxy)propyl]
C-octyl ester, monosodium salt, polymer with butyl 2-propenoate,
methyl 2-methyl-2-propenoate, 2-propenoic acid and
.alpha.-sulfo-.omega.-[1-[(dodecyloxy)methyl]-2-(2propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA
INDEX NAME)

CM 1

CRN 113377-37-0

CMF (C2 H4 O)n C18 H36 O6 S . H3 N

CCI PMS



● NH3

CM 7

CRN 123-34-2 CMF C6 H12 O3

CM 8

CRN 111-87-5 CMF C8 H18 O

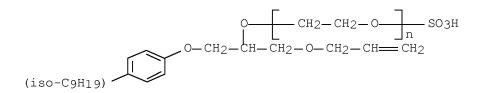
HO- (CH2) 7-Me

RN 354583-17-8 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with 2-hydroxy-3-(2-propenyloxy)propyl octyl hydrogen phosphate monosodium salt and .alpha.-sulfo-.omega.-[1-[(4-isononylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 354583-12-3 CMF (C2 H4 O)n C21 H34 O6 S . H3 N CCI IDS, PMS



● инз

CRN 354552-67-3 CMF C14 H29 O6 P . Na

● Na

CM 3

CRN 141-32-2 CMF C7 H12 O2

RN 354583-18-9 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with 2-hydroxy-3-(2-propenyloxy)propyl octyl hydrogen phosphate monosodium salt and .alpha.-sulfo-.omega.-[4-isononyl(1-propenyl)phenoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 354583-10-1

CMF (C2 H4 O)n C18 H28 O4 S . H3 N

CCI IDS, PMS

$$D1-CH$$
 — $CH-Me$

● инз

CRN 354552-67-3 CMF C14 H29 O6 P . Na

● Na

CM 3

CRN 141-32-2 CMF C7 H12 O2

0 || n-BuO—C— CH**==** CH2

IPCR B01F0017-00 [I,A]; B01F0017-52 [I,A]; C08F0002-16 [I,A]; C08F0002-24
 [I,A]; C08F0002-26 [I,A]; C08F0004-00 [I,A]; C08F0012-04 [I,A];
 C08F0020-12 [I,A]; C08F0216-12 [I,A]; C08F0220-04 [I,A]; C08F0222-16
 [I,A]

CC 37-3 (Plastics Manufacture and Processing)

IT 354552-58-2P 354552-59-3P 354552-61-7P 354552-62-8P

354552-66-2P 354552-68-4P 354583-09-8P 354583-11-2P 354583-13-4P 354583-14-5P 354583-15-6P 354583-16-7P 354583-17-8P 354583-18-9P

(reactive surfactant compns. with controlled CMC for polymer

emulsions)

OS.CITING REF COUNT: 2 THERE ARE 2 CAPLUS RECORDS THAT CITE THIS RECORD (3 CITINGS)

L39 ANSWER 10 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2001:289987 HCAPLUS Full-text

DOCUMENT NUMBER: 134:281635

TITLE: Hydrophilic allylic crosslinking agents for

water-absorbing polymers

INVENTOR(S): Nakamura, Shinichiro; Shimizu, Yasumi; Matsutomi,

Toru

PATENT ASSIGNEE(S): Daiso Co., Ltd., Japan

SOURCE: Jpn. Kokai Tokkyo Koho, 10 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent

10/596.747

LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE		
JP 2001114840	А	20010424	JP 1999-299551	19991021		
			<			
PRIORITY APPLN. INFO.:			JP 1999-299551	19991021		

ED Entered STN: 24 Apr 2001

The agents have 2 allyl groups and .gtoreq.1 hydrophilic group chosen from OSO3Y, AΒ SO3Y, and OP(O)(OY)2(Y = H, alkali metal, ammonium, alk. earth metal). Thus, pentaerythritol triallyl ether (I) was treated with HSO3Cl followed by NaOH to give I sulfate Na salt, which showed 2.21 w/v% soly. in acrylic acid. Acrylic acid was polymd. with I sulfate Na salt to give a polymer showing 46 g/g water absorption.

ΙT 333718-49-3P 333718-50-6P

> (hydrophilic allylic crosslinking agents for water-absorbing polymers)

RN 333718-49-3 HCAPLUS

1-Propanol, 3-(2-propen-1-yloxy)-2,2-bis[(2-propen-1-yloxy)methyl]-, CN 1-(hydrogen sulfate), sodium salt (1:1) (CA INDEX NAME)

$$CH_2-OSO_3H$$
 H_2C — CH_2-O-CH_2-CH — CH_2 — CH_2-CH — CH_2 — C

Na

333718-50-6 HCAPLUS RN

1-Propanol, 3-(2-propenyloxy)-2,2-bis[(2-propenyloxy)methyl]-, CN dihydrogen phosphate, disodium salt (9CI) (CA INDEX NAME)

$$CH_2 - OPO_3H_2$$
 $H_2C = CH - CH_2 - O - CH_2 - C$

●2 Na

333718-51-7P 333718-52-89 333718-53-99 ΙT (hydrophilic allylic crosslinking agents for water-absorbing polymers)

RN 333718-51-7 HCAPLUS CN 2-Propenoic acid, polymer with sodium
3-(2-propenyloxy)-2,2-bis[(2-propenyloxy)methyl]propyl sulfate (9CI)
(CA INDEX NAME)

CM 1

CRN 333718-49-3 CMF C14 H24 O7 S . Na

$$\begin{array}{c} \text{CH}_2\text{--}\,\text{OSO}_3\text{H} \\ \text{H}_2\text{C} = \text{CH}-\text{CH}_2-\text{O}-\text{CH}_2-\text{C}+\text{E}-\text{CH}_2-\text$$

● Na

CM 2

CRN 79-10-7 CMF C3 H4 O2

RN 333718-52-8 HCAPLUS

CN 2-Propenoic acid, polymer with disodium
3-(2-propenyloxy)-2,2-bis[(2-propenyloxy)methyl]propyl phosphate (9CI)
(CA INDEX NAME)

CM 1

CRN 333718-50-6

CMF C14 H25 O7 P . 2 Na

●2 Na

CM 2

CRN 79-10-7 CMF C3 H4 O2

RN 333718-53-9 HCAPLUS

CN 2-Propenoic acid, polymer with sodium 3-(2-propenyloxy)-2,2-bis[(2-propenyloxy)methyl]propyl sulfate, sodium salt (9CI) (CA INDEX NAME)

CM 1

CRN 333718-51-7

CMF (C14 H24 O7 S . C3 H4 O2 . Na)x

CCI PMS

CM 2

CRN 333718-49-3

CMF C14 H24 O7 S . Na

$$\begin{array}{c} \text{CH}_2\text{--}\,\text{OSO}_3\text{H} \\ \text{H}_2\text{C} = \text{CH}-\text{CH}_2-\text{O}-\text{CH}_2-\text{C}+\text{E}-\text{CH}_2-\text$$

● Na

CM 3

CRN 79-10-7 CMF C3 H4 O2

10/596,747

IPCR C08J0003-24 [I,A]; C08F0002-44 [I,A]; C08F0228-02 [I,A]; C08F0230-02
 [I,A]; C08F0291-00 [I,A]; C08L0101-14 [I,A]

CC 37-6 (Plastics Manufacture and Processing)

Section cross-reference(s): 23

IT 333718-49-3P 333718-50-6P

(hydrophilic allylic crosslinking agents for water-absorbing polymers)

IT 333718-51-7P 333718-52-8P 333718-53-9P

(hydrophilic allylic crosslinking agents for water-absorbing polymers)

OS.CITING REF COUNT: 1 THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD (1 CITINGS)

L39 ANSWER 11 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1997:533603 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 127:191884

ORIGINAL REFERENCE NO.: 127:37207a,37210a

TITLE: Compounds and surfactants for emulsifiers for

polymerization and fiber finishing

A DDT TCAMTON NO

בוש עבו

INVENTOR(S): Komiya, Kaoru; Kawamata, Hiromasa; Umezawa, Shohei

PATENT ASSIGNEE(S): Asahi Denka Kogyo K.K., Japan

SOURCE: PCT Int. Appl., 35 pp.

CODEN: PIXXD2

KIND DAME

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

DAMENIM NIO

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
	A1	19970807	₩O 1997-JP186 <	19970128
W: JP, KR, US RW: DE, FR, GB,		10000005	FD 1007 000701	10070120
EP 825167	AI	19980225	EP 1997-900791 <	19970128
EP 825167 R: DE, FR, GB,		20011004		
JP 4031831	B2	20080109	JP 1997-527481 <	19970128
US 5929290	А	19990727	US 1997-913856 <	19970929
JP 2008024942	A	20080207	JP 2007-203331 <	20070803
JP 4824646	В2	20111130		
PRIORITY APPLN. INFO.:			JP 1996-14441 <	A 19960130
			JP 1996-23908 <	A 19960209
			JP 1997-527481 <	A3 19970128
			WO 1997-JP186 <	W 19970128

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

ED Entered STN: 21 Aug 1997

Reactive surfactants RfO(AO)mCH2CH[O(AO')m'X]CH2OCH2CR1:CH2 are prepd., where R1 = H or Me, Rf is a hydrocarbon group or acyl in which .gtoreq.1 H is replaced by F, AO and AO' represent groups selected among C2-4 oxyalkylenes or styrene oxide residues; m, m' = 0 or 1-1,000; and X = H or a hydrophilic group. Thus, C6F13OCH2CH[O(CH2CH2O)30H]CH2OCH2CH:CH2 was prepd. and used as an emulsifier for the polymn. of acrylonitrile, butadiene, and styrene.

IT 194295-78-8P 194295-91-5P 194296-01-0P

(antifogging and antistatic polymers)

RN 194295-78-8 HCAPLUS

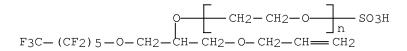
CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2-[(tridecafluorohexyl)oxy]ethoxy]-, potassium salt, polymer with 1-propene, graft (9CI) (CA INDEX NAME)

CM 1

CRN 194294-93-4

CMF (C2 H4 O)n C12 H11 F13 O6 S . K

CCI PMS



• K

CM 2

CRN 115-07-1 CMF C3 H6

H3C−СН**=** СН2

RN 194295-91-5 HCAPLUS

CN Poly(oxy-1,2-ethanediy1), .alpha.-sulfo-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9-hexadecafluoronony1)oxy]methy1]-2(2-propenyloxy)ethoxy]-.omega.-hydroxy-, ammonium salt, polymer with
1-propene, graft (9CI) (CA INDEX NAME)

CM 1

CRN 194295-11-9

CMF (C2 H4 O)n C15 H14 F16 O6 S . H3 N

CCI PMS

■ NH3

CM 2

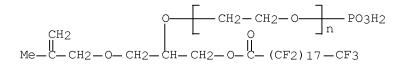
CRN 115-07-1 CMF C3 H6

H3C-CH-CH2

RN 194296-01-0 HCAPLUS
CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15,16,16,17,17,18,18,19,19-heptatriacontafluoro-1oxononadecyl)oxy]methyl]-2-[(2-methyl-2-propenyl)oxy]ethoxy]-,
disodium salt, polymer with 1-propene, graft (9CI) (CA INDEX NAME)

CM 1

CRN 194294-99-0 CMF (C2 H4 O)n C26 H14 F37 O7 P . 2 Na CCI PMS



●2 Na

CM 2

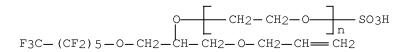
CRN 115-07-1 CMF C3 H6 H3C-CH-CH2

IT 194296-31-6P 194296-39-4P 194296-45-2P (antistatic and antifogging polymers)
RN 194296-31-6 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2-[(tridecafluorohexyl)oxy]ethoxy]-, potassium salt, polymer with ethenylbenzene, graft (9CI) (CA INDEX NAME)

CM 1

CRN 194294-93-4 CMF (C2 H4 O)n C12 H11 F13 O6 S . K CCI PMS



• K

CM 2

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

CCI PMS

RN 194296-39-4 HCAPLUS
CN Poly(oxy-1,2-ethanediy1), .alpha.-sulfo-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9-hexadecafluoronony1)oxy]methy1]-2(2-propenyloxy)ethoxy]-, ammonium salt, polymer with ethenylbenzene, graft (9CI) (CA INDEX NAME)

CM 1

CRN 194295-11-9
CMF (C2 H4 O)n C15 H14 F16 O6 S . H3 N

■ NH3

CM 2

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 194296-45-2 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15,16,16,17,17,18,18,19,19,19-heptatriacontafluoro-1oxononadecyl)oxy]methyl]-2-[(2-methyl-2-propenyl)oxy]ethoxy]-,
disodium salt, polymer with ethenylbenzene, graft (9CI) (CA INDEX NAME)

CM 1

CRN 194294-99-0

CMF (C2 H4 O)n C26 H14 F37 O7 P . 2 Na $\,$

CCI PMS

●2 Na

CM 2

CRN 100-42-5 CMF C8 H8 H_2C \longrightarrow CH \longrightarrow Ph

IT 194296-09-8P 194296-18-9P 194296-25-8P (fibers; antisoiling fabrics)

RN 194296-09-8 HCAPLUS

CN 1,4-Benzenedicarboxylic acid, polymer with 1,2-ethanediol and .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2[(tridecafluorohexyl)oxy]ethoxy]poly(oxy-1,2-ethanediyl) potassium salt, graft (9CI) (CA INDEX NAME)

CM 1

CRN 194294-93-4 CMF (C2 H4 O)n C12 H11 F13 O6 S . K CCI PMS

• K

CM 2

CRN 107-21-1 CMF C2 H6 O2

HO-CH2-CH2-ОН

CM 3

CRN 100-21-0 CMF C8 H6 O4

RN 194296-18-9 HCAPLUS

CN 1,4-Benzenedicarboxylic acid, polymer with 1,2-ethanediol and .alpha.-sulfo-.omega.-[1-[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9-hexadecafluorononyl)oxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt, graft (9CI) (CA INDEX NAME)

CM 1

CRN 194295-11-9

CMF (C2 H4 O)n C15 H14 F16 O6 S . H3 N

CCI PMS

● NH3

CM 2

CRN 107-21-1 CMF C2 H6 O2

HO-CH2-CH2-OH

CM 3

CRN 100-21-0 CMF C8 H6 O4

RN 194296-25-8 HCAPLUS

CN 1,4-Benzenedicarboxylic acid, polymer with 1,2-ethanediol and .alpha.-phosphono-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15,16,16,17,17,18,18,19,19,19-heptatriacontafluoro-1oxononadecyl)oxy]methyl]-2-[(1-methyl-2-propenyl)oxy]ethoxy]poly(oxy-

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1,2-ethanediyl) disodium salt, graft (9CI) (CA INDEX NAME)

CM 1

CRN 194294-99-0

CMF (C2 H4 O)n C26 H14 F37 O7 P . 2 Na

CCI PMS

●2 Na

CM 2

CRN 107-21-1 CMF C2 H6 O2

 $HO-CH_2-CH_2-OH$

CM 3

CRN 100-21-0 CMF C8 H6 O4

IT 194295-16-4P 194295-20-0P 194295-23-3P

194295-26-6P 194295-30-2P 194295-33-5P

194295-36-8P 194295-40-4P 194295-43-7P

(manuf. and reactive surfactants as emulsifiers for)

RN 194295-16-4 HCAPLUS

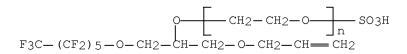
CN 2-Propenenitrile, polymer with 1,3-butadiene, ethenylbenzene and .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2-

[(tridecafluorohexyl)oxy]ethoxy]poly(oxy-1,2-ethanediyl) potassium
salt (9CI) (CA INDEX NAME)

CRN 194294-93-4

CMF (C2 H4 O)n C12 H11 F13 O6 S . K

CCI PMS



K

CM 2

CRN 107-13-1 CMF C3 H3 N

 $H2C \longrightarrow CH - C \longrightarrow N$

CM 3

CRN 106-99-0 CMF C4 H6

H2C== CH- CH== CH2

CM 4

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 194295-20-0 HCAPLUS

CN 2-Propenenitrile, polymer with 1,3-butadiene, ethenylbenzene and .alpha.-sulfo-.omega.-[1-[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9-

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hexadecafluorononyl)oxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 194295-11-9

CMF (C2 H4 O)n C15 H14 F16 O6 S . H3 N $\,$

CCI PMS

● NH3

CM 2

CRN 107-13-1

CMF C3 H3 N

$$H_2C \longrightarrow CH - C \longrightarrow N$$

CM 3

CRN 106-99-0 CMF C4 H6

 H_2C \longrightarrow CH \longrightarrow CH_2

CM 4

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

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RN 194295-23-3 HCAPLUS

CM 1

CRN 194294-99-0

CMF (C2 H4 O)n C26 H14 F37 O7 P . 2 Na

CCI PMS

•2 Na

CM 2

CRN 107-13-1 CMF C3 H3 N

H2C==CH-C==N

CM 3

CRN 106-99-0 CMF C4 H6

H2C== CH- CH== CH2

CM 4

CRN 100-42-5 CMF C8 H8 H2C=CH-Ph

RN 194295-26-6 HCAPLUS
CN 2-Propenoic acid, 2-methyl-, methyl ester, polymer with
.alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2[(tridecafluorohexyl)oxy]ethoxy]poly(oxy-1,2-ethanediyl) potassium
salt (9CI) (CA INDEX NAME)

CM 1

CRN 194294-93-4
CMF (C2 H4 O)n C12 H11 F13 O6 S . K
CCI PMS

● K

CM 2

CRN 80-62-6

CMF C5 H8 O2

H2C O || || || || |Me-C-C-OMe

CCI PMS

RN 194295-30-2 HCAPLUS
CN 2-Propenoic acid, 2-methyl-, methyl ester, polymer with
.alpha.-sulfo-.omega.-[1-[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9-hexadecafluorononyl)oxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 194295-11-9
CMF (C2 H4 O)n C15 H14 F16 O6 S . H3 N

■ NH3

CM 2

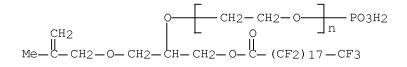
CRN 80-62-6 CMF C5 H8 O2

RN 194295-33-5 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, methyl ester, polymer with .alpha.-phosphono-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15,16,16,17,17,18,18,19,19,19-heptatriacontafluoro-1oxononadecyl)oxy]methyl]-2-[(2-methyl-2-propenyl)oxy]ethoxy]poly(oxy-1,2-ethanediyl) disodium salt (9CI) (CA INDEX NAME)

CM 1

CRN 194294-99-0 CMF (C2 H4 O)n C26 H14 F37 O7 P . 2 Na CCI PMS



•2 Na

CM 2

CRN 80-62-6 CMF C5 H8 O2

RN 194295-36-8 HCAPLUS

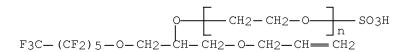
CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2-[(tridecafluorohexyl)oxy]ethoxy]-, potassium salt, polymer with chloroethene (9CI) (CA INDEX NAME)

CM 1

CRN 194294-93-4

CMF (C2 H4 O)n C12 H11 F13 O6 S . K

CCI PMS



• K

CM 2

CRN 75-01-4 CMF C2 H3 C1

H2C=CH-C1

RN 194295-40-4 HCAPLUS

CN Poly(oxy-1,2-ethanediy1), .alpha.-sulfo-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9-hexadecafluoronony1)oxy]methy1]-2(2-propenyloxy)ethoxy]-, ammonium salt, polymer with chloroethene
(9CI) (CA INDEX NAME)

CM 1

CRN 194295-11-9

CMF (C2 H4 O)n C15 H14 F16 O6 S . H3 N

CCI PMS

■ NH3

CM 2

CRN 75-01-4 CMF C2 H3 C1

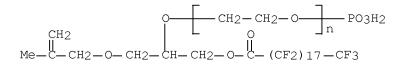
H2C==CH-C1

RN 194295-43-7 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15,16,16,17,17,18,18,19,19,19-heptatriacontafluoro-1oxononadecyl)oxy]methyl]-2-[(2-methyl-2-propenyl)oxy]ethoxy]-,
disodium salt, polymer with chloroethene (9CI) (CA INDEX NAME)

CM 1

CRN 194294-99-0 CMF (C2 H4 O)n C26 H14 F37 O7 P . 2 Na CCI PMS



●2 Na

CM 2

CRN 75-01-4 CMF C2 H3 C1 H2C=CH-C1

194295-46-0P 194295-50-6P 194295-53-9P ΙT 194295-56-2P 194295-64-2P 194295-70-0P (manuf. of antifogging and antistatic polymers and reactive surfactants as emulsifiers for) RN 194295-46-0 HCAPLUS Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(2-CN propenyloxy)methyl]-2-[(tridecafluorohexyl)oxy]ethoxy]-, potassium salt, polymer with ethenylbenzene (9CI) (CA INDEX NAME) CM 1 CRN 194294-93-4 CMF (C2 H4 O)n C12 H11 F13 O6 S . K CCI PMS

• K

CM 2

CRN 100-42-5

CMF C8 H8

H2C=CH-Ph

RN 194295-50-6 HCAPLUS
CN Poly(oxy-1,2-ethanediy1), .alpha.-sulfo-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9-hexadecafluoronony1)oxy]methy1]-2(2-propenyloxy)ethoxy]-, ammonium salt, polymer with ethenylbenzene
(9CI) (CA INDEX NAME)

CM 1

CRN 194295-11-9
CMF (C2 H4 O)n C15 H14 F16 O6 S . H3 N
CCI PMS

■ NH3

CM 2

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 194295-53-9 HCAPLUS
CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15,16,16,17,17,18,18,19,19,19-heptatriacontafluoro-1oxononadecyl)oxy]methyl]-2-[(2-methyl-2-propenyl)oxy]ethoxy]-,

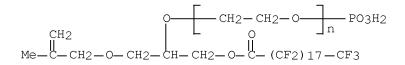
oxononadecy1)oxy]methy1]-2-[(2-methy1-2-propeny1)oxy]ethoxy]-, disodium salt, polymer with ethenylbenzene (9CI) (CA INDEX NAME)

CM 1

CRN 194294-99-0

CMF (C2 H4 O)n C26 H14 F37 O7 P . 2 Na

CCI PMS



●2 Na

CM 2

CRN 100-42-5

CMF C8 H8

 H_2C \longrightarrow CH \longrightarrow Ph

RN 194295-56-2 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, polymer with butyl 2-propenoate and .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2[(tridecafluorohexyl)oxy]ethoxy]poly(oxy-1,2-ethanediyl) potassium salt (9CI) (CA INDEX NAME)

CM 1

CRN 194294-93-4

CMF (C2 H4 O)n C12 H11 F13 O6 S . K

CCI PMS

• K

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 79-41-4 CMF C4 H6 O2

RN 194295-64-2 HCAPLUS

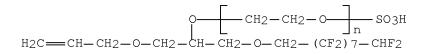
CN 2-Propenoic acid, 2-methyl-, polymer with butyl 2-propenoate and .alpha.-sulfo-.omega.-[1-[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9-hexadecafluorononyl)oxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 194295-11-9

CMF (C2 H4 O)n C15 H14 F16 O6 S . H3 N

CCI PMS



● инз

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 79-41-4 CMF C4 H6 O2

RN 194295-70-0 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, polymer with butyl 2-propenoate and .alpha.-phosphono-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15,16,16,17,17,18,18,19,19,19-heptatriacontafluoro-1oxononadecyl)oxy]methyl]-2-[(1-methyl-2-propenyl)oxy]ethoxy]poly(oxy-1,2-ethanediyl) disodium salt (9CI) (CA INDEX NAME)

CRN 194294-99-0

CMF (C2 H4 O)n C26 H14 F37 O7 P . 2 Na

CCI PMS

●2 Na

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 79-41-4 CMF C4 H6 O2

IT 194294-90-1P 194294-91-2P 194294-92-3P 194294-93-4P 194294-94-5P 194294-95-6P 194294-96-7P 194294-97-8P 194294-98-9P 194294-99-0P 194295-01-7P

(reactive surfactants for emulsifiers for polymn. and fiber finishing)

RN 194294-90-1 HCAPLUS

194295-11-9P

CN 2-Propanol, 1-(2-propen-1-yloxy)-3-[(1,1,2,2,3,3,4,4,5,5,6,6,6-tridecafluorohexyl)oxy]-, 2-(hydrogen sulfate), sodium salt (1:1) (CA INDEX NAME)

Na

RN 194294-91-2 HCAPLUS

CN Heptanoic acid, 2,2,3,3,4,4,5,5,6,6,7,7,7-tridecafluoro-, 3-(2-propen-1-yloxy)-2-(sulfooxy)propyl ester, sodium salt (1:1) (CA INDEX NAME)

Na

RN 194294-92-3 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2-[(tridecafluorohexyl)oxy]ethoxy]- (9CI) (CA INDEX NAME)

RN 194294-93-4 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2-[(tridecafluorohexyl)oxy]ethoxy]-, potassium salt (9CI) (CA INDEX NAME)

• K

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RN 194294-94-5 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2-[(2,2,3,3,4,4,5,5,6,6,7,7,7-tridecafluoro-1-oxoheptyl)oxy]ethoxy]- (9CI) (CA INDEX NAME)

RN 194294-95-6 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(2-propenyloxy)methyl]-2-[(2,2,3,3,4,4,5,5,6,6,7,7,7-tridecafluoro-1-oxoheptyl)oxy]ethoxy]-, potassium salt (9CI) (CA INDEX NAME)

$$c_{13}c_{-1}$$
 c_{-1} c_{-

• K

RN 194294-96-7 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1-[[(heptatriacontafluorooctadecyl)oxy]methyl]-2-[(2-methyl-2propenyl)oxy]ethoxy]-, disodium salt (9CI) (CA INDEX NAME)

●2 Na

RN 194294-97-8 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1-[[(heptatriacontafluorooctadecyl)oxy]methyl]-2-[(2-methyl-2propenyl)oxy]ethoxy]-, diammonium salt (9CI) (CA INDEX NAME)

●2 NH3

RN 194294-98-9 HCAPLUS

CN Poly(oxy-1,2-ethanediy1), .alpha.-phosphono-.omega.-[1-[[(heptatriacontafluorooctadecy1)oxy]methy1]-2-[(2-methy1-2propeny1)oxy]ethoxy]-, magnesium salt (9CI) (CA INDEX NAME)

Mg

RN 194294-99-0 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15
,16,16,17,17,18,18,19,19,19-heptatriacontafluoro-1oxononadecyl)oxy]methyl]-2-[(2-methyl-2-propenyl)oxy]ethoxy]-,
disodium salt (9CI) (CA INDEX NAME)

●2 Na

RN 194295-00-6 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15,16,16,17,17,18,18,19,19-heptatriacontafluoro-1oxononadecyl)oxy]methyl]-2-[(2-methyl-2-propenyl)oxy]ethoxy]-,
diammonium salt (9CI) (CA INDEX NAME)

●2 NH3

RN 194295-01-7 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15
,16,16,17,17,18,18,19,19,19-heptatriacontafluoro-1oxononadecyl)oxy]methyl]-2-[(2-methyl-2-propenyl)oxy]ethoxy]-,
magnesium salt (9CI) (CA INDEX NAME)

Ma

RN 194295-11-9 HCAPLUS

CN Poly(oxy-1,2-ethanediy1), .alpha.-sulfo-.omega.-[1-[[(2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9-hexadecafluoronony1)oxy]methy1]-2-(2-propenyloxy)ethoxy]-, ammonium salt (9CI) (CA INDEX NAME)

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194295-87-9P
                194295-91-5P
                               194295-94-8P 194295-97-1P
    194296-01-0P 194296-05-4P
       (antifogging and antistatic polymers)
ΙT
    194296-29-2P 194296-31-6P
                              194296-33-8P
                                           194296-35-0P
    194296-41-8P 194296-43-0P
    194296-45-2P 194296-47-4P
       (antistatic and antifogging polymers)
    194296-07-6P 194296-09-8P 194296-11-2P 194296-13-4P
ΤТ
    194296-15-6P 194296-18-9P
                              194296-20-3P 194296-22-5P
    194296-25-8P 194296-27-0P
       (fibers; antisoiling fabrics)
ΙT
    194295-15-3P 194295-16-4P 194295-17-5P 194295-18-6P
    194295-23-3P 194295-24-4P 194295-25-5P
    194295-26-6P 194295-27-7P 194295-28-8P 194295-29-9P
    194295-30-2P 194295-31-3P 194295-32-4P
    194295-33-5P 194295-34-6P
                               194295-35-7P
    194295-36-8P 194295-37-9P 194295-38-0P 194295-39-1P
    194295-40-4P 194295-41-5P 194295-42-6P
    194295-43-7P 194295-44-8P
       (manuf. and reactive surfactants as emulsifiers for)
    ΤТ
    194295-49-3P 194295-50-6P 194295-51-7P
                                           194295-52-8P
                194295-54-0P
                              194295-55-1P
    194295-53-9P
    194295-56-2P 194295-58-4P 194295-60-8P 194295-62-0P
    194295-64-2P 194295-66-4P 194295-68-6P
    194295-70-0P 194295-72-2P
       (manuf. of antifogging and antistatic polymers and reactive
       surfactants as emulsifiers for)
ΙT
    194294-82-1P 194294-83-2P 194294-84-3P
                                             194294-85-4P
    194294-86-5P
                 194294-87-6P
                               194294-88-7P
                                           194294-89-8P
    194294-90-1P 194294-91-2P 194294-92-3P
    194294-93-4P 194294-94-5P 194294-95-6P
    194294-96-7P 194294-97-8P 194294-98-9P
    194294-99-0P 194295-00-6P 194295-01-7P
    194295-02-8P 194295-03-9P 194295-04-0P
                                           194295-05-1P
                194295-07-3P 194295-08-4P
    194295-06-2P
                                            194295-09-5P
    194295-10-8P
                 194295-11-9P
                              194295-12-0P
                                            194295-13-1P
    194295-14-2P
                194368-99-5P 194369-00-1P 194369-01-2P
    194369-02-3P
       (reactive surfactants for emulsifiers for polymn. and fiber
       finishing)
OS.CITING REF COUNT:
                      1
                           THERE ARE 1 CAPLUS RECORDS THAT CITE THIS
                           RECORD (1 CITINGS)
                           THERE ARE 2 CITED REFERENCES AVAILABLE FOR
REFERENCE COUNT:
                      2
                           THIS RECORD. ALL CITATIONS AVAILABLE IN THE
                           RE FORMAT
L39 ANSWER 12 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN
ACCESSION NUMBER:
                      1996:248322 HCAPLUS Full-text
DOCUMENT NUMBER:
                      124:345042
ORIGINAL REFERENCE NO.: 124:64091a,64094a
TITLE:
                      Storage-stable aqueous polymer compositions for
                      coating films with good water and chemical
                      resistances and strength
INVENTOR(S):
                      Nakahata, Takashi; Nakada, Tadahiro; Oka, Masashi
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10/596.747

PATENT ASSIGNEE(S): Asahi Denka Kogyo KK, Japan SOURCE: Jpn. Kokai Tokkyo Koho, 12 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 08034897	А	19960206	JP 1994-172225	19940725
			<	
JP 3508030	В2	20040322		
PRIORITY APPLN. INFO.:			JP 1994-172225	19940725
			/	

ED Entered STN: 27 Apr 1996

AB Title compns. (solids content 2-90%) contain (A) 1-85% polyurethanes and (B) 1-85% acrylic polymers obtained by polymg. acrylic unsatd. monomer mixts. in the presence of reactive emulsifiers having copolymerizable unsatd. bonds. The reactive emulsifiers may be CH2:CR1XCH2CH(OZ)CH2O(AO)mR2 [R1 = H, Me; R2 = C6-30 hydrocarbyl, acyl; A = C2-4 alkylene; X = CH2O, C(O)O, CH2OC(O); Z = H, nonionic or anionic hydrophilic group; m = 0-100]. Thus, 200 parts aq. polyurethane (prepd. from polypropylene glycol 49, dicyclohexylmethane diisocyanate 176, dimethylolpropionic acid 70, N-methylpyrrolidone 196, Et3N 48, hexamethylenediamine 5, and H2O 456 parts) and 233 parts acrylic emulsion [prepd. from Me methacrylate 45, Bu acrylate 45, glycidyl methacrylate 10, CH2:CHCH2OCH2CH[O(C2H4O)10SO3NH4]CH2OC6H4C9H19-p 3, (NH4)2S2O8 0.6, and H2O 130 parts] were mixed to give a 40% storage-stable aq. compn., which was applied on a glass plate to give a coating film with good water and chem. resistances.

IT 176744-68-6 176744-69-7 176744-70-0 176744-71-1 176744-75-5 176744-77-7

(storage-stable aq. polyurethane-acrylic polymer compns. for coating films with good water and chem. resistances and strength)

RN 176744-68-6 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, methyl ester, polymer with butyl 2-propenoate, oxiranylmethyl 2-methyl-2-propenoate and .alpha.-sulfo-.omega.-[1-[(4-nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 136931-77-6

CMF (C2 H4 O)n C21 H34 O6 S . H3 N

CCI PMS

● NH3

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 106-91-2 CMF C7 H10 O3

CM 4

CRN 80-62-6 CMF C5 H8 O2

RN 176744-69-7 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, methyl ester, polymer with butyl 2-propenoate, ethenylbenzene, N-(hydroxymethyl)-2-propenamide, oxiranylmethyl 2-methyl-2-propenoate and .alpha.-sulfo-.omega.-[1-[(4-nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 136931-77-6 CMF (C2 H4 O)n C21 H34 O6 S . H3 N CCI PMS

● NH3

CM 2

CRN 924-42-5 CMF C4 H7 N O2

CM 3

CRN 141-32-2 CMF C7 H12 O2

CM 4

CRN 106-91-2 CMF C7 H10 O3

CM 5

CRN 100-42-5 CMF C8 H8

CRN 80-62-6 CMF C5 H8 O2

RN 176744-70-0 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, methyl ester, polymer with ethenylbenzene, 2-ethylhexyl 2-propenoate,
N-(hydroxymethyl)-2-propenamide, oxiranylmethyl 2-methyl-2-propenoate and .alpha.-sulfo-.omega.-[1-[(4-nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 136931-77-6

CMF (C2 H4 O)n C21 H34 O6 S . H3 N

CCI PMS

● NH3

CM 2

CRN 924-42-5 CMF C4 H7 N O2

CRN 106-91-2 CMF C7 H10 O3

$$\overset{\text{O}}{\longleftarrow}_{\text{CH}_2-\text{O}-}\overset{\text{O}}{\underset{\text{CH}_2}{\parallel}}\overset{\text{CH}_2}{\underset{\text{CH}_2}{\parallel}}$$

CM 4

CRN 103-11-7 CMF C11 H20 O2

$$CH_2 - O - C - CH = CH_2$$

 $Et - CH - Bu - n$

CM 5

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

CM 6

CRN 80-62-6 CMF C5 H8 O2

RN 176744-71-1 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, methyl ester, polymer with butyl 2-propenoate and .alpha.-sulfo-.omega.-[1-[(4-nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 136931-77-6

CMF (C2 H4 O)n C21 H34 O6 S . H3 N

CCI PMS

$$HO_3S$$
 $-CH_2-CH_2$ $-CH_2-CH_2-CH_2-CH_2$ $-CH_2-CH_2-CH_2-CH_2$

● NH3

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 80-62-6 CMF C5 H8 O2

RN 176744-75-5 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, methyl ester, polymer with butyl 2-propenoate, oxiranylmethyl 2-methyl-2-propenoate and .alpha.-phosphono-.omega.-[1-[(dodecyloxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) diammonium salt (9CI) (CA INDEX NAME)

CRN 176744-74-4

CMF (C2 H4 O)n C18 H37 O6 P . 2 H3 N

CCI PMS

●2 NH3

CM 2

CRN 141-32-2

CMF C7 H12 O2

CM 3

CRN 106-91-2 CMF C7 H10 O3

CM 4

CRN 80-62-6 CMF C5 H8 O2

RN 176744-77-7 HCAPLUS

CN Butanedioic acid, sulfo-, 1-dodecyl 4-(2-propenyl) ester, sodium salt, polymer with butyl 2-propenoate, methyl 2-methyl-2-propenoate and oxiranylmethyl 2-methyl-2-propenoate (9CI) (CA INDEX NAME)

CM 1

CRN 162880-23-1

CMF C19 H34 O7 S . Na

● Na

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 106-91-2 CMF C7 H10 O3

CM 4

CRN 80-62-6 CMF C5 H8 O2



IPCI C08L0033-14 [ICM,6]; C08F0002-24 [ICS,6]; C08L0075-04 [ICS,6]

IPCR C08L0033-14 [I,A]; C08F0002-24 [I,A]; C08L0033-04 [I,A]; C08L0075-00

[I,A]; C08L0075-04 [I,A]

CC 37-6 (Plastics Manufacture and Processing)

Section cross-reference(s): 38, 42

IT 103336-45-4 113988-50-4 143482-33-1 157609-04-6 176744-65-3

176744-66-4 176744-67-5 176744-68-6 176744-69-7 176744-70-0 176744-71-1 176744-73-3 176744-75-5 176744-76-6

(storage-stable aq. polyurethane-acrylic polymer compns. for coating films with good water and chem. resistances and strength)

L39 ANSWER 13 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1995:654989 HCAPLUS Full-text

DOCUMENT NUMBER: 123:35414

ORIGINAL REFERENCE NO.: 123:6490h,6491a

TITLE: Heat blocking-resistant coatings INVENTOR(S): Ikebayashi, Nobuhiko; Koshio, Takeaki

PATENT ASSIGNEE(S): Hoechst Gosei KK, Japan

Ι

SOURCE: Jpn. Kokai Tokkyo Koho, 10 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

176744-77-7

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 07026165	А	19950127	JP 1993-220382	19930706
			<	
PRIORITY APPLN. INFO.:			JP 1993-220382	19930706

ED Entered STN: 05 Jul 1995

GΙ

AB Title coatings contain .alpha., .beta.-ethylenic monomer-based polymers prepd. in the presence of emulsifiers CH2:CR1CH2OCH2CH(OZ)CH2O(AO)mR2 (A = C2-4 alkylene;

10/596,747

R1 = H, Me; R2 = C8-24 hydrocarbyl or acyl; Z = H, nonionic or anionic hydrophilic groups), pigments, and colloidal SiO2 with Y .gtoreq.[25 - (1/4 .times. X)] and X = 5-90, in which X wt.% = pigments/(SiO2 composite polymers + pigments) and Y wt.% = SiO2/[SiO2 composite polymers - (polymers + SiO2)]. A mixt. of 13-nm SiO2 particles and an emulsion contg. 100-nm polymer particles prepd. from Bu acrylate, Me methacrylate, and methacrylic acid in the presence of I was further mixed with TiO2, CaCO3, and other additives gave a compn. with X = 60% and Y = 29.4% and showing good heat blocking resistance and storage stability at 50.degree. for 1 wk.

IT 164463-57-4 164463-58-5 164463-59-6

164463-60-9 164463-62-1

(silica composite; coatings with heat blocking resistance and storage stability)

RN 164463-57-4 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, polymer with butyl 2-propenoate, methyl 2-methyl-2-propenoate and .alpha.-sulfo-.omega.-[1-[[4-(nonyloxy)phenoxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 164463-56-3

CMF (C2 H4 O)n C21 H34 O7 S . H3 N

CCI PMS

● NH3

CM 2

CRN 141-32-2 CMF C7 H12 O2

0 || n-BuO-C-CH-CH2

CM 3

CRN 80-62-6 CMF C5 H8 O2

CRN 79-41-4 CMF C4 H6 O2

RN 164463-58-5 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, polymer with butyl 2-propenoate, ethenyltriethoxysilane, methyl 2-methyl-2-propenoate and .alpha.-sulfo-.omega.-[1-[[4-(nonyloxy)phenoxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 164463-56-3

CMF (C2 H4 O)n C21 H34 O7 S . H3 N $\,$

CCI PMS

● инз

CM 2

CRN 141-32-2 CMF C7 H12 O2

CRN 80-62-6 CMF C5 H8 O2

CM 4

CRN 79-41-4 CMF C4 H6 O2

CM 5

CRN 78-08-0 CMF C8 H18 O3 Si

RN 164463-59-6 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, polymer with butyl 2-propenoate, ethenylbenzene, methyl 2-methyl-2-propenoate and .alpha.-sulfo-.omega.-[1-[[4-(nonyloxy)phenoxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

10/596,747

CRN 164463-56-3 CMF (C2 H4 O)n C21 H34 O7 S . H3 N CCI PMS

● инз

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

CM 4

CRN 80-62-6 CMF C5 H8 O2

CM 5

CRN 79-41-4 CMF C4 H6 O2

RN 164463-60-9 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, polymer with butyl 2-propenoate, methyl 2-methyl-2-propenoate, 2-propenoic acid and .alpha.-sulfo-.omega.-[1-[[4-(nonyloxy)phenoxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 164463-56-3

CMF (C2 H4 O)n C21 H34 O7 S . H3 N

CCI PMS

● инз

CM 2

CRN 141-32-2 CMF C7 H12 O2

0 || n-BuO-C-CH-CH2 CM 3

CRN 80-62-6 CMF C5 H8 O2

CM 4

CRN 79-41-4 CMF C4 H6 O2

CM 5

CRN 79-10-7 CMF C3 H4 O2

RN 164463-62-1 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, polymer with butyl 2-propenoate, methyl 2-methyl-2-propenoate and .alpha.-phosphono-.omega.-[1-[[4-(nonyloxy)phenoxy]methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) disodium salt (9CI) (CA INDEX NAME)

CM 1

CRN 164463-61-0

CMF (C2 H4 O)n C21 H35 O7 P . 2 Na

CCI PMS

●2 Na

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 80-62-6 CMF C5 H8 O2

CM 4

CRN 79-41-4 CMF C4 H6 O2

CC 42-7 (Coatings, Inks, and Related Products)

IT 164463-57-4 164463-58-5 164463-59-6

164463-60-9 164463-62-1

(silica composite; coatings with heat blocking resistance and storage stability)

L39 ANSWER 14 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1995:571394 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 123:229313

ORIGINAL REFERENCE NO.: 123:40977a,40980a

TITLE: Dispersing agents for suspension polymerization of

vinyl chloride monomers

INVENTOR(S): Mizutari, Takeaki; Tsuzuki, Masahide; Komya, Kaoru

PATENT ASSIGNEE(S): Asahi Denka Kogyo KK, Japan SOURCE: Jpn. Kokai Tokkyo Koho, 7 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 07062005	А	19950307	JP 1993-211809	19930826
			<	
PRIORITY APPLN. INFO.:			JP 1993-211809	19930826
			/	

ED Entered STN: 25 May 1995

Title agents which do not remain in the product polymers in a free state comprise C2:CR1CH2OCH2CH(OX)CH2O(AO)nR2 [sic; A = C2-4 alkylene; R1 = H, Me; R2 = C1-24 hydrocarbyl, acyl; n = 0-50; X = hydrophilic group of (AO)mH, (AO)rSO3M, or (AO)kP(O)(OM1)(OM2); M1, M2 = H, alkali metal, alk. earth metal, (org.) ammonium; m = 1-100; r, k = 1-50]. Thus, equimolar nonylphenol and allyl glycidyl ether were reacted at 90 .+-. 5.degree. for 5 h and then 1 mol the resulting product was further reacted with 10 mol ethylene oxide to obtain a dispersing agent (I). Then 100 parts vinyl chloride was polymd. in H2O in the presence of 1 part I and di-2-ethylhexyl peroxydicarbonate at 57.degree. for 7 h showing no scale deposition on the reactor wall. The resulting polymer showed good water resistance and thermal stability.

IT 111144-58-2P 168009-71-0P

(reactive dispersants for manuf. of thermally stable water-resistant vinyl chloride polymers)

RN 111144-58-2 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1[(nonylphenoxy)methyl]-2-(2-propen-1-yloxy)ethoxy]-, sodium salt (1:1)
(CA INDEX NAME)



 $D1 - (CH_2) 8 - Me$

● Na

RN 168009-71-0 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]-, sodium salt (9CI) (CA INDEX NAME)



D1 - (CH₂) 8 - Me

●x Na

IT 168009-72-1P 168009-73-2P 168009-74-3P 168009-75-4P 168009-76-5P 168109-70-4P

168109-71-5P 169970-88-1P

(reactive dispersants for manuf. of thermally stable water-resistant vinyl chloride polymers)

RN 168009-72-1 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]-, sodium salt, polymer with chloroethene (9CI) (CA INDEX NAME)

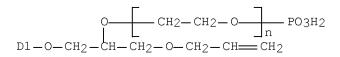
CM 1

CRN 168009-71-0

CMF (C2 H4 O)n C21 H35 O6 P . x Na



D1 - (CH2) 8 - Me



●x Na

CM 2

CRN 75-01-4 CMF C2 H3 C1

H2C=CH-C1

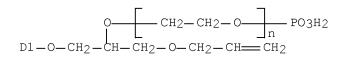
RN 168009-73-2 HCAPLUS
CN Poly(oxy-1,2-ethanediy1), .alpha.-phosphono-.omega.-[1[(nonylphenoxy)methy1]-2-(2-propenyloxy)ethoxy]-, polymer with chloroethene and ethene (9CI) (CA INDEX NAME)

CM 1

CRN 168009-71-0 CMF (C2 H4 O)n C21 H35 O6 P . x Na CCI IDS, PMS



D1 - (CH2) 8 - Me



●x Na

CM 2

CRN 75-01-4 CMF C2 H3 C1

H2C=CH-C1

CM 3

CRN 74-85-1 CMF C2 H4

H2C=CH2

RN 168009-74-3 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1- [(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]-, sodium salt, polymer with chloroethene and ethene (9CI) (CA INDEX NAME)

CM 1

CRN 111144-58-2

CMF (C2 H4 O)n C21 H34 O6 S . Na



$$D1 - (CH_2) 8 - Me$$

● Na

CM 2

CRN 75-01-4 CMF C2 H3 C1

H2C=CH-C1

CM 3

CRN 74-85-1 CMF C2 H4

H2C=CH2

RN 168009-75-4 HCAPLUS

CN Acetic acid ethenyl ester, polymer with chloroethene and .alpha.-phosphono-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) sodium salt (9CI) (CA INDEX NAME)

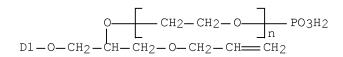
CM 1

CRN 168009-71-0

CMF (C2 H4 O)n C21 H35 O6 P . x Na



D1 - (CH2) 8 - Me



●x Na

CM 2

CRN 108-05-4 CMF C4 H6 O2

ACO-CH-CH2

CM 3

CRN 75-01-4 CMF C2 H3 C1

H2C==CH-C1

RN 168009-76-5 HCAPLUS

CN Acetic acid ethenyl ester, polymer with chloroethene and .alpha.-sulfo-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy][poly(oxy-1,2-ethanediyl)] sodium salt (9CI) (CA INDEX NAME)

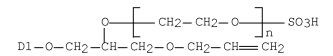
CM 1

CRN 111144-58-2

CMF (C2 H4 O)n C21 H34 O6 S . Na



$$D1 - (CH_2) 8 - Me$$



● Na

CM 2

CRN 108-05-4 CMF C4 H6 O2

ACO-CH-CH2

CM 3

CRN 75-01-4 CMF C2 H3 C1

H2C==CH-C1

RN 168109-70-4 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]-, sodium salt, polymer with chloroethene and ethoxyethene (9CI) (CA INDEX NAME)

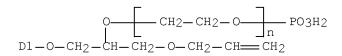
CM 1

CRN 168009-71-0

CMF (C2 H4 O)n C21 H35 O6 P . x Na



D1 - (CH2) 8 - Me



●x Na

CM 2

CRN 109-92-2 CMF C4 H8 O

H3C-CH2-O-CH-CH2

CM 3

CRN 75-01-4 CMF C2 H3 C1

H2C==CH-C1

RN 168109-71-5 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1- [(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]-, sodium salt, polymer with chloroethene and ethoxyethene (9CI) (CA INDEX NAME)

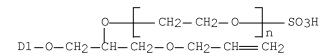
CM 1

CRN 111144-58-2

CMF (C2 H4 O)n C21 H34 O6 S . Na



$$D1 - (CH_2) 8 - Me$$



● Na

CM 2

CRN 109-92-2 CMF C4 H8 O

H3C-CH2-O-CH CH2

CM 3

CRN 75-01-4 CMF C2 H3 C1

H2C==CH-C1

RN 169970-88-1 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]-, sodium salt, polymer with chloroethene (9CI) (CA INDEX NAME)

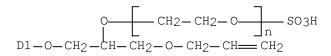
CM 1

CRN 111144-58-2

CMF (C2 H4 O)n C21 H34 O6 S . Na



D1 - (CH2) 8 - Me



Na

CM 2

CRN 75-01-4 CMF C2 H3 C1

H2C== CH- C1

ACCESSION NUMBER:

```
IPCI C08F0002-18 [ICM,6]; B01F0017-42 [ICS,6]; C08F0014-06 [ICS,6]
IPCR B01F0017-42 [I,A]; C08F0002-18 [I,A]; C08F0014-00 [I,A]; C08F0014-06
     [I,A]
    35-4 (Chemistry of Synthetic High Polymers)
CC
    Section cross-reference(s): 46
ΙT
    75-21-8DP, Oxirane, reaction products with glycerol (meth)allyl ethers
    75-56-9DP, reaction products with glycerol (meth)allyl ethers
    26249-20-7DP, Butylene oxide, reaction products with glycerol
     (meth)allyl ethers 111144-58-2P
                                       111144-60-6P
    168009-71-0P
                   168111-42-0P
                                 168399-97-1P
        (reactive dispersants for manuf. of thermally stable
       water-resistant vinyl chloride polymers)
    75-01-4DP, polymers with alkoxylated glycerol (meth)allyl ethers
ΤТ
    79-10-7DP, 2-Propenoic acid, esters, polymers with vinyl chloride and
    alkoxylated glycerol (meth)allyl ethers 168009-69-6P 168009-70-9P
    168009-72-1P
                   168009-73-2P
                                  168009-74-3P
    168009-75-4P
                   168009-76-5P
                                  168036-63-3P
    168109-69-1P 168109-70-4P
                                  168109-71-5P
                                  168253-58-5P
    168111-44-2P
                  168112-60-5P
                                                168397-31-7P
    168397-32-8P
                  168397-33-9P
                                  168609-02-7P
                                                169970-88-1P
    199542-29-5P
        (reactive dispersants for manuf. of thermally stable
       water-resistant vinyl chloride polymers)
```

L39 ANSWER 15 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN

1993:192325 HCAPLUS Full-text

DOCUMENT NUMBER: 118:192325

ORIGINAL REFERENCE NO.: 118:33065a,33068a

TITLE: Synthesis of polymerizable surfactant and its

application to emulsion polymerization

AUTHOR(S): Yokota, Kinya; Ichihara, Akinobu; Shinike, Hitoshi

CORPORATE SOURCE: Appl. Lab., Dai-Ichi Kogyo Seiyaku Co. Ltd.,

Kyoto, 600, Japan

SOURCE: Special Publication - Royal Society of Chemistry

(1992), 107(Ind. Appl. Surfactants III), 29-48

CODEN: SROCDO; ISSN: 0260-6291

DOCUMENT TYPE: Journal LANGUAGE: English ED Entered STN: 14 May 1993

AB As polymerizable surfactants, ethoxylated

1-nonylphenoxy-2-hydroxy-3-(allyloxy) propane, its sulfate, and its phosphate were synthesized. Ethoxylated 2-allyl-4-nonylphenol, its sulfate, and its phosphate were also synthesized. The yields of these polymerizable surfactants were >95%. These polymerizable surfactants were applied as emulsifiers for emulsion polymn. of Et acrylate and Bu acrylate/styrene. Their polymn. stabilities were the same as those of surfactants commonly used and great improvements were seen in low foaming of the emulsion and water resistance of the polymer film.

IT 136931-77-6P 146847-16-7P 146847-17-8P

(prepn. and surfactant properties of polymerizable)

RN 136931-77-6 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(4-nonylphenoxy)methyl]-2-(2-propen-1-yloxy)ethoxy]-, ammonium salt (1:1) (CA INDEX NAME)

$$\begin{array}{c|c} \text{HO}_{3}\text{S} & \hline & \text{O-CH}_2\text{-CH}_2 \\ \hline & \text{H}_2\text{C} = \text{CH-CH}_2\text{-O-CH}_2\text{-CH-CH}_2\text{-O} \\ \end{array}$$

● NH3

RN 146847-16-7 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1-[(4-nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]- (9CI) (CA INDEX NAME)

RN 146847-17-8 HCAPLUS

CN Poly(oxy-1,2-ethanediy1), .alpha.,.alpha.'-phosphinicobis[.omega.-[1-[(4-nonylphenoxy)methy1]-2-(2-propen-1-yloxy)ethoxy]- (CA INDEX NAME)

PAGE 1-B

$$-CH_2-CH_2$$
 $-CH_2-CH_2-CH_2$ $-CH_2-CH_2-CH_2$ $-CH_2-CH_2-CH_2$ $-CH_2-CH_2-CH_2-CH_2$ $-CH_2-CH_2-CH_2-CH_2-CH_2$

IT 146847-26-9P 146847-31-6P (prepn. of)

RN 146847-26-9 HCAPLUS

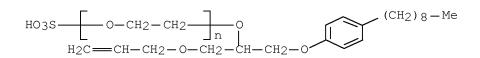
CN 2-Propenoic acid, ethyl ester, polymer with .alpha.-sulfo-.omega.-[1-[(4-nonylphenoxy)methyl]-2-(2-propen-1-yloxy)ethoxy][poly(oxy-1,2-ethanediyl)] ammonium salt (1:1) (CA INDEX NAME)

CM 1

CRN 136931-77-6

CMF (C2 H4 O)n C21 H34 O6 S . H3 N

CCI PMS



● инз

CM 2

CRN 140-88-5 CMF C5 H8 O2

RN 146847-31-6 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with .alpha.-sulfo-.omega.-[1-[(4-nonylphenoxy)methyl]-2-(2-propen-1-yloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (1:1) (CA INDEX NAME)

CM 1

CRN 136931-77-6

CMF (C2 H4 O)n C21 H34 O6 S . H3 N

CCI PMS

● NH3

CM 2

CRN 141-32-2 CMF C7 H12 O2

CC 35-2 (Chemistry of Synthetic High Polymers)

Section cross-reference(s): 46

IT 136931-77-6P 140651-97-4P 146847-16-7P 146847-17-8P 146847-18-9P 146847-19-0P

(prepn. and surfactant properties of polymerizable)

IT 140651-98-5P 146847-26-9P 146847-28-1P 146847-29-2P

(prepn. of)

OS.CITING REF COUNT: 5 THERE ARE 5 CAPLUS RECORDS THAT CITE THIS RECORD (5 CITINGS)

L39 ANSWER 16 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1992:450024 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 117:50024

ORIGINAL REFERENCE NO.: 117:8937a,8940a

TITLE: Oriented polyester films containing acrylic

resin-treated inorganic particles

INVENTOR(S): Kuze, Katsuro; Matsumoto, Haruo; Murashige,

Ryuichi

PATENT ASSIGNEE(S): Toyobo Co., Ltd., Japan; Nippon Magphane Co., Ltd.

SOURCE: Jpn. Kokai Tokkyo Koho, 9 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
 JР 04039336	 A	19920210	JP 1990-145041	19900601
01 04037330	Α	19920210	<	19900001
JP 2767485	В2	19980618	TD 1000 145041	10000601
PRIORITY APPLN. INFO.:			JP 1990-145041	19900601

ED Entered STN: 08 Aug 1992

The title films contain 0.005-3.0% inert inorg. particles having improved compatibility as a result of treatment with copolymers of .gtoreq.1 monomer H2C:CR1CO2X (R1 = H, Me, HOCH2CH2; X = H, monovalent or divalent metal, ammonium, amine) and .gtoreq.1 allyl ether H2C:CHCH2OCH2[CH[CH2(OR3)rZ]]p(OR2)qY [R2-3 = C2-4 alkylene; Y, Z = OH, C1-4 alkoxy, [monovalent or divalent metal-, ammonium salt-, org. amine (salt)-, or C1-4 alkyl ester-substituted] monovalent phosphate group or sulfate group; Y and Z may form divalent phosphate, sulfonate, or sulfate group]. A mixt. of CaCO3 and 98:2 acrylic

acid-H2C:CHCH2OCH2CH[(OCH2CH2)nOPO3H2]CH2(OCH2CH2)nOPO3H2 copolymer was spray dried to prep. a filler, and a polyester was prepd. from terephthalic acid and ethylene glycol in the presence of the filler, giving a product which formed a film having dynamic friction coeff. 0.38, haze 5.7%, and void ratio 0.25%.

IT 142551-11-9 142551-13-1 142551-82-4 142571-36-6

(fillers modified by, for compatibility in oriented polyester films)

RN 142551-11-9 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, sodium salt, polymer with 1-[(2-propenyloxy)methyl]-1,2-ethanediyl bis(hydrogen sulfate) (9CI) (CA INDEX NAME)

CM 1

CRN 142551-10-8 CMF C6 H12 O9 S2

OSO3H HO3SO-CH2-CH-CH2-O-CH2-CH=CH2

CM 2

CRN 5536-61-8 CMF C4 H6 O2 . Na

● Na

RN 142551-13-1 HCAPLUS
CN 2-Propenoic acid, ammonium salt, polymer with
 .alpha.,.alpha.'-[1-[(2-propenyloxy)methyl]-1,2-ethanediyl]bis[.omega. (sulfooxy)poly(oxy-1,2-ethanediyl)] sodium salt (9CI) (CA INDEX NAME)

CM 1

CRN 142551-12-0

CMF (C2 H4 O)n (C2 H4 O)n C6 H12 O9 S2 . x Na
CCI PMS

H2C=CH-CH2-O-CH2
H03SO
$$\begin{array}{c} CH_2 - CH_2 -$$

CM 2

CRN 10604-69-0 CMF C3 H4 O2 . H3 N

● NH3

RN 142551-82-4 HCAPLUS
CN 2-Propenoic acid, polymer with
.alpha.,.alpha.'-[1-[(2-propenyloxy)methyl]-1,2-ethanediyl]bis[.omega.-

(phosphonooxy)poly(oxy-1,2-ethanediyl)] (9CI) (CA INDEX NAME)

CM 1

CRN 90717-19-4

CMF (C2 H4 O)n (C2 H4 O)n C6 H14 O9 P2

CCI PMS

CM 2

CRN 79-10-7 CMF C3 H4 O2

RN 142571-36-6 HCAPLUS

CN 2-Propenoic acid, sodium salt, polymer with 2-hydroxy-3-(2-propenyloxy)-1-propanesulfonic acid monopotassium salt (9CI) (CA INDEX NAME)

CM 1

CRN 84019-66-9 CMF C6 H12 O5 S . K

• ĸ

CM 2

CRN 7446-81-3

CMF C3 H4 O2 . Na

0 HO-C-CH=CH2

Na

IPCR C08J0005-18 [I,A]; B29C0055-02 [I,A]; B29K0067-00 [N,A]; B29L0007-00
 [N,A]; C08F0016-14 [I,A]; C08F0016-16 [I,A]; C08F0216-14 [I,A];
 C08F0216-16 [I,A]; C08K0009-04 [I,A]; C08L0067-00 [I,A]; C08L0067-02
 [I,A]

CC 37-6 (Plastics Manufacture and Processing)

Section cross-reference(s): 38

IT 142551-11-9 142551-13-1 142551-82-4

142551-83-5 142571-36-6

(fillers modified by, for compatibility in oriented polyester films)

L39 ANSWER 17 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1992:428378 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 117:28378

ORIGINAL REFERENCE NO.: 117:5123a,5126a

TITLE: Allyl ether copolymers as coupling agents for

inert fillers in oriented polyester films for

improving sliding properties

INVENTOR(S): Kuze, Katsuro; Matsumoto, Haruo; Murashige,

Ryuichi

PATENT ASSIGNEE(S): Toyobo Co., Ltd., Japan; Nippon Magphane Co., Ltd.

SOURCE: Jpn. Kokai Tokkyo Koho, 8 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 04033933	А	19920205	JP 1990-140667	19900529
			<	
PRIORITY APPLN. INFO.:			JP 1990-140667	19900529
			<	

ED Entered STN: 26 Jul 1992

The title couplers are obtained from copolymers of allyl ethers CH2:CHCH2OCH2(ZH)p(OZ1)qR1 [I; Z = C[CH2(OZ2)rR2]; p = 1-4; q, r = 0-100; Z1-2 = C2-4 alkylene; R1-2 = OH, alkoxy, monophosphoric acid, salts or esters, and monosulfonic acid, salts or esters, or R1-2 together can form a divalent similar phosphate and sulfonate group in the above], (meth)acrylic acids, their salts or esters, and unsatd. dicarboxylic acids selected from maleic, fumaric or itaconic acids, their salts or diesters. Thus, mixing an aq. soln. of a 80:18:2 acrylic acid-diammonium maleate-3-allyloxy-1,2- di(polyoxyethylene)propanephosphate (i.e. I, with q, r = .apprx.4) copolymer with CaCO3 in solids wt. ratio 0.005:1, and spray drying gave treated particles. Biaxially oriented PET film contq. 0.25%

of the particles had surface roughness 0.025 .mu.m, dynamic friction coeff. 0.37 .mu.d, haze 5.0%, void ratio 0.17%, and no. of broken void 0.05/mm2, compared to 0.020, 0.43, 13.2, 1.35, and 1.2, resp. for similar film contg. fillers without the I treatment.

IT 142357-64-0 142357-65-1 142357-67-3

(couplers, for inert fillers in polyester films for tapes, for improved sliding properties)

RN 142357-64-0 HCAPLUS

CN 2-Butenedioic acid (2Z)-, diammonium salt, polymer with 2-propenoic acid and .alpha.,.alpha.'-[1-[(2-propenyloxy)methyl]-1,2-ethanediyl]bis[.omega.-(phosphonooxy)poly(oxy-1,2-ethanediyl)] (9CI) (CA INDEX NAME)

CM 1

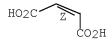
CRN 90717-19-4 CMF (C2 H4 O)n (C2 H4 O)n C6 H14 O9 P2 CCI PMS

$$H_2C = CH - CH_2 - O - CH_2$$
 $H_2O_3PO = CH_2 - CH_2 - O - CH_2 - CH_2 - CH_2 - O - CH_2 -$

CM 2

CRN 23705-99-9 CMF C4 H4 O4 . 2 H3 N

Double bond geometry as shown.



●2 NH3

CM 3

CRN 79-10-7 CMF C3 H4 O2

RN 142357-65-1 HCAPLUS

CN 2-Butenedioic acid (2E)-, disodium salt, polymer with 2-hydroxy-3-(2-propenyloxy)-1-propanesulfonic acid monosodium salt and sodium 2-propenoate (9CI) (CA INDEX NAME)

CM 1

CRN 52556-42-0 CMF C6 H12 O5 S . Na

Na

CM 2

CRN 17013-01-3 CMF C4 H4 O4 . 2 Na

Double bond geometry as shown.

●2 Na

CM 3

CRN 7446-81-3 CMF C3 H4 O2 . Na

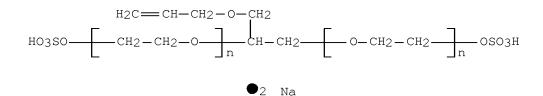
● Na

RN 142357-67-3 HCAPLUS

CN 2-Butenedioic acid (2Z)-, disodium salt, polymer with 2-propenoic acid and .alpha.,.alpha.'-[1-[(2-propenyloxy)methyl]-1,2-ethanediyl]bis[.omega.-(sulfooxy)poly(oxy-1,2-ethanediyl)] disodium salt (9CI) (CA INDEX NAME)

CM 1

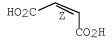
CRN 142357-66-2 CMF (C2 H4 O)n (C2 H4 O)n C6 H12 O9 S2 . 2 Na CCI PMS



CM 2

CRN 371-47-1 CMF C4 H4 O4 . 2 Na

Double bond geometry as shown.



•2 Na

CM 3

CRN 79-10-7 CMF C3 H4 O2

IPCI C08J0005-18 [ICM,5]; C08K0009-04 [ICS,5]; C08L0067-02 [ICS,5]; C08L0067-00 [ICI,5]

IPCR C08J0005-18 [I,A]; C08K0009-04 [I,A]; C08L0067-00 [I,A]; C08L0067-02 [I,A]

CC 38-3 (Plastics Fabrication and Uses)

142357-64-0 142357-65-1 142357-67-3 ΙT

> (couplers, for inert fillers in polyester films for tapes, for improved sliding properties)

L39 ANSWER 18 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1988:133883 HCAPLUS <u>Full-text</u>
DOCUMENT NUMBER: 108:133883

ORIGINAL REFERENCE NO.: 108:21953a,21956a

Surface-active compounds having a polymerizable TITLE:

moiety from allyl glycidyl ether

Yokota, Kinya; Ichihara, Akinobu INVENTOR(S):

Daiichi Kogyo Seiyaku Co., Ltd., Japan PATENT ASSIGNEE(S):

Eur. Pat. Appl., 19 pp. SOURCE:

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 244841	A2	19871111	EP 1987-106533	19870506
EP 244841	А3	19890607		
EP 244841 R: DE, FR, GB	В1	19920318		
JP 63054927	А	19880309	JP 1986-105119 <	19860507
JP 63054928	А	19880309	JP 1986-118956 <	19860522
JP 63054929	A	19880309	JP 1986-121954 <	19860526
JP 62279833	A	19871204	JP 1986-124305	19860528
JP 63080837	A	19880411	JP 1986-127006 <	19860530
JP 02022695	В	19900521		
US 4814514	A	19890321	US 1987-46353 <	19870506
JP 63126535	А	19880530	JP 1987-198778 <	19870807
JP 05076335	В	19931022		
JP 63214336	А	19880907	JP 1987-198779 <	19870807
JP 04068013	В	19921030		
JP 63199271	A	19880817	JP 1987-208689 <	19870821
JP 2589502	В2	19970312		
JP 63203872	А	19880823	JP 1987-208690 <	19870821
JP 04002714	В	19920120		

JP 63205384	А	19880824	JP 1987-208691 <		19870821
JP 05015753	В	19930302	`		
US 4939283	A	19900703	US 1988-284588		19881215
			<		
PRIORITY APPLN. INFO.:			JP 1986-105119	Α	19860507
			<		
			JP 1986-118956	А	19860522
			<		
			JP 1986-121954	A	19860526
			<		
			JP 1986-124305	Α	19860528
			<		
			JP 1986-127006	А	19860530
			<		
			US 1987-46353	А3	19870506
			<		

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

ED Entered STN: 15 Apr 1988

AB Surfactants R1O(AO)xCH2CH[O(AO)yH]CH2OCH2CR2:CH2 (R1 = alkyl, alkenyl, alkyl aryl, or arylakylaryl contg. 8-30 C; R2 = H or Me; A = C2-4 alkylene; x = 0-100; y = 1-200) and their sulfate, phosphate, and sulfosuccinate esters are prepd. The surfactants are useful as emulsifiers in the emulsion or suspension polymn. of ethylenically unsatd. monomers, as finishing agents for hydrophobic textiles, as antistatic agents for plastics, as pigment dispersants, etc. A reaction product of 1.0 mol nonylphenol and 1.0 mol allyl glycidyl ether was ethoxylated with 10 mol oxirane and used as a copolymerizable emulsifier in the polymn. of a Bu acrylate-styrene mixt., giving a stable emulsion which was dried and cured at 110.degree. to give a water-resistant coating.

IT 113356-42-6P 113356-44-8P 113356-45-9P

113377-36-9P 113377-37-0P

113377-38-1P 113377-63-2P 113405-84-8P

113405-85-9P 113405-86-0P 113405-87-1P

113441-08-0P

(prepn. and surface activity of)

RN 113356-42-6 HCAPLUS

CN Oxirane, methyl-, polymer with oxirane, bis(1-phenylethyl)phenyl 2-hydroxy-3-(2-propenyloxy)propyl ether, diether with .alpha.,.alpha.'-(1,4-dioxo-2-sulfo-1,4-butanediyl)bis[.omega.-hydroxypoly(oxy-1,2-ethanediyl)], sodium salt, block (9CI) (CA INDEX NAME)

CM 1

CRN 177072-56-9

CMF (C2 H4 O)n (C2 H4 O)n C16 H26 O11 S

CCI PMS

PAGE 1-B

$$- \begin{picture}(20,2) \put(0,0){\line(1,0){100}} \put(0,0){\line(1,0){1$$

CM 2

CRN 25640-70-4 CMF C22 H22 O

CCI IDS

1/2 (D1-OH)

CM 3

CRN 106392-12-5 CMF (C3 H6 O . C2 H4 O)x CCI PMS

CM 4

CRN 75-56-9 CMF C3 H6 O



CM 5

CRN 75-21-8 CMF C2 H4 O



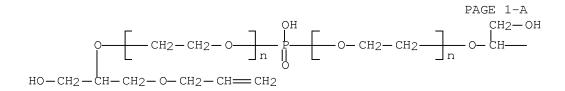
RN 113356-44-8 HCAPLUS
CN Oxirane, methyl-, polymer with oxirane, bis(1-phenylethyl)phenyl 2-hydroxy-3-(2-propenyloxy)propyl ether, diether with .alpha.,.alpha.'-phosphinicobis[.omega.-hydroxypoly(oxy-1,2-ethanediyl)], block (9CI) (CA INDEX NAME)

CM 1

CRN 177072-59-2

CMF (C2 H4 O)n (C2 H4 O)n C12 H23 O8 P

CCI PMS



PAGE 1-B

— CH2-O-CH2-CH= CH2

CM 2

CRN 25640-70-4 CMF C22 H22 O CCI IDS

1/2 (D1-OH)

CM 3

CRN 106392-12-5

CMF (C3 H6 O . C2 H4 O)x

CCI PMS

CM 4

CRN 75-56-9 CMF C3 H6 O



CM 5

CRN 75-21-8 CMF C2 H4 O



RN 113356-45-9 HCAPLUS

CN Oxirane, methyl-, polymer with oxirane, bis(1-phenylethyl)phenyl 2-hydroxy-3-(2-propenyloxy)propyl ether, ether with .alpha.-sulfo-.omega.-hydroxypoly(oxy-1,2-ethanediyl), ammonium salt, block (9CI) (CA INDEX NAME)

CM 1

$$1/2 (D1-OH)$$



CM 5

CRN 75-21-8 CMF C2 H4 O



RN 113377-36-9 HCAPLUS

CN Poly(oxy-1,2-ethanediy1), .alpha.-phosphono-.omega.-[1- [(dodecyloxy)methy1]-2-(2-propenyloxy)ethoxy]- (9CI) (CA INDEX NAME)

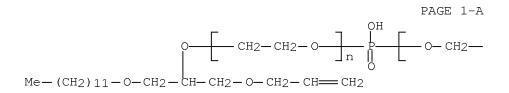
RN 113377-37-0 HCAPLUS

CN Poly(oxy-1,2-ethanediy1), .alpha.-sulfo-.omega.-[1- [(dodecyloxy)methy1]-2-(2-propenyloxy)ethoxy]-, ammonium salt (9CI) (CA INDEX NAME)

● NH3

RN 113377-38-1 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.,.alpha.'-phosphinicobis[.omega.-[1-[(dodecyloxy)methyl]-2-(2-propenyloxy)ethoxy]- (9CI) (CA INDEX NAME)



$$-CH_2$$
 $-CH_2$ $-CH_2$ $-CH_2$ $-CH_2$ $-CH_2$ $-CH_2$ $-CH_3$ $-CH_4$ $-CH_4$ $-CH_5$ $-CH_5$ $-CH_6$ $-CH_$

RN 113377-63-2 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.,.alpha.'-(1,4-dioxo-2-sulfo-1,4-butanediyl)bis[.omega.-[1-[(dodecyloxy)methyl]-2-(2-propenyloxy)ethoxy]-, sodium salt (9CI) (CA INDEX NAME)

● Na

RN 113405-84-8 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-phosphono-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]- (9CI) (CA INDEX NAME)



$$D1 - (CH_2) 8 - Me$$

$$O-CH_2-CH_2-O-CH_2-O-CH_2-CH=CH_2$$

RN 113405-85-9 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propen-1-yloxy)ethoxy]-, ammonium salt (1:1) (CA INDEX NAME)



$$D1 - (CH_2) 8 - Me$$

● NH3

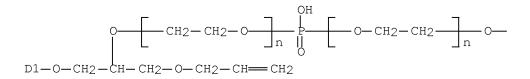
RN 113405-86-0 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.,.alpha.'-phosphinicobis[.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]- (9CI) (CA INDEX NAME)

PAGE 1-A

$$2\left[\bigcirc \right]$$

$$2 \left[D1-(CH_2)8-Me \right]$$



PAGE 1-B

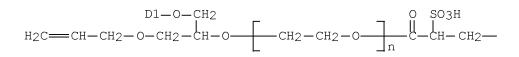
RN 113405-87-1 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.,.alpha.'-(1,4-dioxo-2-sulfo-1,4-butanediyl)bis[.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]-, sodium salt (9CI) (CA INDEX NAME)

PAGE 1-A



 $2 \Gamma D1-(CH_2)_8-Me$



● Na

PAGE 1-B

$$-\frac{0}{C} - \frac{CH_2 - O - D1}{CH - CH_2 - O - CH_2 - CH_2} - \frac{CH_2 - O - D1}{CH - CH_2 - O - CH_2 - CH_2} - CH_2$$

RN 113441-08-0 HCAPLUS

CN Oxirane, methyl-, polymer with oxirane, bis(1-phenylethyl)phenyl 2-hydroxy-3-(2-propenyloxy)propyl ether, ether with .alpha.-phosphono-.omega.-hydroxypoly(oxy-1,2-ethanediyl) (9CI) (CA INDEX NAME)

CM 1

CRN 177072-61-6

CMF (C2 H4 O)n C6 H13 O6 P

CCI PMS

CM 2

CRN 25640-70-4

CMF C22 H22 O

CCI IDS

1/2 (D1-OH)

CM 3

CRN 9003-11-6

CMF (C3 H6 O . C2 H4 O) x

CCI PMS

CM 4

CRN 75-56-9

CMF C3 H6 O



CM 5

CRN 75-21-8 CMF C2 H4 O



CM 1

CRN 113377-37-0

CMF (C2 H4 O)n C18 H36 O6 S . H3 N

CCI PMS

О——— CH2— CH2— О——— SO3H

■ NH3

CM 2

CRN 108-05-4 CMF C4 H6 O2

AcO-CH-CH2

RN 113377-49-4 HCAPLUS
CN Acetic acid ethenyl ester, polymer with

.alpha.,.alpha.'-phosphinicobis[.omega.-[1-[(dodecyloxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl)] and

.alpha.-phosphono-.omega.-[1-[(dodecyloxy)methyl]-2-(2propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) (9CI) (CA INDEX NAME)

CM 1

CRN 113377-38-1

CMF (C2 H4 O)n (C2 H4 O)n C36 H71 O8 P

CCI PMS

PAGE 1-A

OH

CH2-CH2-O-In

OH

O-CH2
Me-(CH2)11-O-CH2-CH-CH2-O-CH2-CH=CH2

PAGE 1-B

$$-CH_2 - O - (CH_2)_{11} - Me$$
 $-CH_2 - CH_2 - CH_2 - CH_2 - CH_2 - CH_2 - CH_2$

CM 2

CRN 113377-36-9

CMF (C2 H4 O)n C18 H37 O6 P

CCI PMS

CM 3

CRN 108-05-4 CMF C4 H6 O2

AcO-CH \longrightarrow CH2

CCI PMS

RN 113377-64-3 HCAPLUS
CN Acetic acid ethenyl ester, polymer with
.alpha.,.alpha.'-(1,4-dioxo-2-sulfo-1,4-butanediyl)bis[.omega.-[1[(dodecyloxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl)]
sodium salt (9CI) (CA INDEX NAME)

CM 1

CRN 113377-63-2
CMF (C2 H4 O)n (C2 H4 O)n C40 H74 O11 S . Na

PAGE 1-B

Na

CH2-O- (CH2)11-Me -CH2-CH2-CH2-O-CH-CH2-CH=-CH2

CM 2

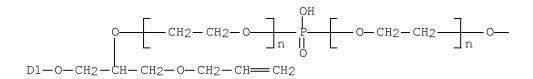
CRN 108-05-4 CMF C4 H6 O2

AcO-CH \longrightarrow CH2

ΙT 113405-89-3P 113405-90-6P 113405-92-8P 113405-93-9P 113405-95-1P 113405-96-2P 113405-97-3P 113405-98-4P 113405-99-5P 113431-94-02 113431-93-99 (prepn. of self-emulsifiable, as coating materials) 113405-89-3 HCAPLUS RN 2-Propenoic acid, ethyl ester, polymer with CN .alpha.,.alpha.'-phosphinicobis[.omega.-[1-[(nonylphenoxy)methyl]-2-(2propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl)] and .alpha.-phosphono-.omega.-[1-[(nonylphenoxy)methyl]-2-(2propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) (9CI) (CA INDEX NAME) СМ 1 CRN 113405-86-0 (C2 H4 O)n (C2 H4 O)n C42 H67 O8 P CMF CCI IDS, PMS

PAGE 1-A





PAGE 1-B

CM 2

CRN 113405-84-8

CMF (C2 H4 O)n C21 H35 O6 P



D1-(CH2)8-Me

CM 3

CRN 140-88-5 CMF C5 H8 O2

RN 113405-90-6 HCAPLUS

CN 2-Propenoic acid, ethyl ester, polymer with .alpha.-sulfo-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 113405-85-9

CMF (C2 H4 O)n C21 H34 O6 S . H3 N

CCI IDS, PMS



 $D1 - (CH_2) 8 - Me$

● NH3

CRN 140-88-5 CMF C5 H8 O2

RN 113405-92-8 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with ethenylbenzene and .alpha.-sulfo-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA INDEX NAME)

CM 1

CRN 113405-85-9

CMF (C2 H4 O)n C21 H34 O6 S . H3 N $\,$

CCI IDS, PMS



$$D1 - (CH_2) 8 - Me$$

● NH3

CM 2

CRN 141-32-2 CMF C7 H12 O2

```
CM 3

CRN 100-42-5

CMF C8 H8
```

H2C=CH-Ph

CM

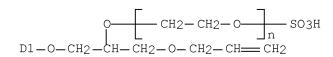
1

RN 113405-93-9 HCAPLUS
CN 2-Propenoic acid, butyl ester, polymer with ethenylbenzene,
 .alpha.-[1-[(nonylphenoxy)ethyl]-2-(2-propenyloxy)ethyl]-.omega. hydroxypoly(oxy-1,2-ethanediyl) and
 .alpha.-sulfo-.omega.-[1-[(nonylphenoxy)methyl]-2-(2 propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA
 INDEX NAME)

CRN 113405-85-9 CMF (C2 H4 O)n C21 H34 O6 S . H3 N CCI IDS, PMS



 $D1 - (CH_2) 8 - Me$



● NH3

CM 2

CRN 111144-60-6

CMF (C2 H4 O)n C21 H34 O3

CCI IDS, PMS



$$H_2C$$
 = $CH - CH_2 - O - CH_2 - CH - CH_2 - CH_2$

CM 3

CRN 141-32-2 CMF C7 H12 O2

CM 4

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 113405-95-1 HCAPLUS
CN 2-Propenoic acid, methyl ester, polymer with
 .alpha.-sulfo-.omega.-[1-[(nonylphenoxy)methyl]-2-(2 propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) ammonium salt (9CI) (CA
 INDEX NAME)

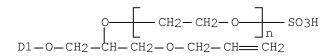
CM 1

CRN 113405-85-9

CMF (C2 H4 O)n C21 H34 O6 S . H3 N $\,$



$$D1 - (CH_2) 8 - Me$$



● NH3

CM 2

CRN 96-33-3 CMF C4 H6 O2

RN 113405-96-2 HCAPLUS

CN 2-Propenoic acid, methyl ester, polymer with

.alpha.,.alpha.'-phosphinicobis[.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl)] and

.alpha.-phosphono-.omega.-[1-[(nonylphenoxy)methyl]-2-(2propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) (9CI) (CA INDEX NAME)

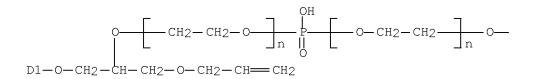
CM 1

CRN 113405-86-0

CMF (C2 H4 O)n (C2 H4 O)n C42 H67 O8 P

PAGE 1-A





PAGE 1-B

$$_{\rm CH_2-O-D1}^{\rm CH_2-O-D1}_{\rm CH-CH_2-O-CH_2-CH=CH_2}^{\rm CH}$$

CM 2

CRN 113405-84-8

CMF (C2 H4 O)n C21 H35 O6 P



D1-(CH2)8-Me

CM 3

CRN 96-33-3 CMF C4 H6 O2

RN 113405-97-3 HCAPLUS
CN 2-Propenoic acid, ethyl ester, polymer with
 .alpha.,.alpha.'-(1,4-dioxo-2-sulfo-1,4-butanediyl)bis[.omega.-[1 [(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2 ethanediyl)] sodium salt (9CI) (CA INDEX NAME)

CM 1

CRN 113405-87-1

CMF (C2 H4 O)n (C2 H4 O)n C46 H70 O11 S . Na

PAGE 1-A

$$2\left[\bigcirc \right]$$

$$2 \Gamma D1-(CH_2)_8-Me$$

Na

PAGE 1-B

$$O = CH_2 - CH_$$

CM 2

CRN 140-88-5 CMF C5 H8 O2

RN 113405-98-4 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with

10/596,747

.alpha.,.alpha.'-(1,4-dioxo-2-sulfo-1,4-butanediyl)bis[.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl)] sodium salt and ethenylbenzene (9CI) (CA INDEX NAME)

CM 1

CRN 113405-87-1 CMF (C2 H4 O)n (C2 H4 O)n C46 H70 O11 S . Na CCI IDS, PMS

PAGE 1-A



$$2 \left[D1 - (CH_2)_8 - Me \right]$$

Na

PAGE 1-B

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 113405-99-5 HCAPLUS

CN 2-Propenoic acid, methyl ester, polymer with .alpha.,.alpha.'-(1,4-dioxo-2-sulfo-1,4-butanediyl)bis[.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl)] sodium salt (9CI) (CA INDEX NAME)

CM 1

CRN 113405-87-1

CMF (C2 H4 O)n (C2 H4 O)n C46 H70 O11 S . Na

CCI IDS, PMS

PAGE 1-A



$$2 \begin{bmatrix} D1-(CH_2)8-Me \end{bmatrix}$$

$$\begin{array}{c} \text{D1-O-CH2} \\ \text{H2C-CH-CH2-O-CH2-CH-O-CH2-CH2-O-J}_n \\ \end{array} \\ \begin{array}{c} \text{O} \\ \text{SO3H} \\ \text{C-CH-CH2-CH2-O-J}_n \\ \end{array}$$

Na

PAGE 1-B

$$- \stackrel{\text{O}}{\text{C}} - \stackrel{\text{C}}{\text{C}} - \text{C} + 2 - \text{C} + 2$$

CM 2

CRN 96-33-3 CMF C4 H6 O2

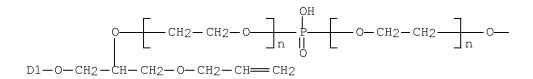
RN 113431-93-9 HCAPLUS

CM 1

CRN 113405-86-0 CMF (C2 H4 O)n (C2 H4 O)n C42 H67 O8 P CCI IDS, PMS

PAGE 1-A





PAGE 1-B

$$_{\rm CH_2-O-D1}^{\rm CH_2-O-D1}_{\rm CH-CH_2-O-CH_2-CH=CH_2}^{\rm CH}$$

CM 2

CRN 113405-84-8

CMF (C2 H4 O)n C21 H35 O6 P



D1-(CH2)8-Me

CM 3

CRN 141-32-2 CMF C7 H12 O2

CM 4

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 113431-94-0 HCAPLUS
CN 2-Propenoic acid, butyl ester, polymer with ethenylbenzene,
 .alpha.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethyl]-.omega. hydroxypoly(oxy-1,2-ethanediyl) and
 .alpha.,alpha.'-phosphinicobis[.omega.-[1-[(nonylphenoxy)methyl]-2-(2 propenyloxy)ethyl]poly(oxy-1,2-ethanediyl)] (9CI) (CA INDEX NAME)

CM 1

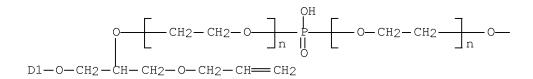
CRN 113405-86-0

CMF (C2 H4 O)n (C2 H4 O)n C42 H67 O8 P

CCI IDS, PMS

PAGE 1-A





PAGE 1-B

$$_{\rm CH_2-O-D1}^{\rm CH_2-O-D1}_{\rm CH-CH_2-O-CH_2-CH=CH_2}^{\rm CH}$$

CM 2

CRN 111144-60-6

CMF (C2 H4 O)n C21 H34 O3



$$D1 - (CH_2) 8 - Me$$

$$D1-O-CH_2$$
 H_2C
 $CH-CH_2-O-CH_2-CH$
 $O-CH_2-CH_2$
 $O-CH_2-CH_2$
 $O-CH_2-CH_2$

CM 3

CRN 141-32-2 CMF C7 H12 O2

CM 4

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

```
IPCI C07C0043-23 [ICM,4]; C07C0043-178 [ICS,4]; C07C0141-08 [ICS,4];
    C07C0143-12 [ICS,4]; C07F0009-09 [ICS,4]; C08G0065-28 [ICS,4];
    C08F0002-24 [ICS, 4]
IPCR B01F0017-00 [I,A]; C07C0043-11 [I,A]; C07C0043-178 [I,A]; C07C0043-23
     [I,A]; C07C0305-06 [I,A]; C07C0305-10 [I,A]; C07C0309-17 [I,A];
    C07F0009-09 [I,A]; C08F0002-26 [I,A]; C08G0065-26 [I,A]; C08G0065-327
    [I,A]; C08G0065-334 [I,A]
CC
    46-4 (Surface Active Agents and Detergents)
    Section cross-reference(s): 35, 37, 38, 40, 42
ΙT
                   111144-60-6P
                                  113356-42-6P
                                                113356-43-7P
    111100-57-3P
                  113356-45-9P
    113356-44-8P
                                  113356-46-0P
    113356-47-1P
                   113356-48-2P
                                  113356-49-3P
                                                113356-50-6P
                                  113377-36-9P
    113356-51-7P
                   113356-52-8P
    113377-37-0P
                  113377-38-1P
                                  113377-63-2P
    113405-81-5P
                  113405-83-7P 113405-84-8P
    113405-85-9P 113405-86-0P
                                  113405-87-1P
```

10/596,747

113441-08-0P 113441-10-4P

113405-88-2P 113431-90-6P

```
(prepn. and surface activity of)
    113377-23-4P 113377-48-3P 113377-49-4P 113377-64-3P 113472-80-3P
ΙT
       (prepn. of self-emulsifiable, as adhesives)
ΙT
    111144-66-2P 111144-70-8P 113405-89-3P
    113405-90-6P 113405-91-7P 113405-92-8P
    113405-93-9P 113405-94-0P 113405-95-1P
    113405-96-2P 113405-97-3P 113405-98-4P
    113405-99-5P 113431-93-9P 113431-94-0P
    113473-94-2P 113473-95-3P 113473-96-4P
        (prepn. of self-emulsifiable, as coating materials)
OS.CITING REF COUNT: 19
                           THERE ARE 19 CAPLUS RECORDS THAT CITE THIS
                             RECORD (20 CITINGS)
L39 ANSWER 19 OF 19 HCAPLUS COPYRIGHT 2012 ACS on STN
ACCESSION NUMBER:
                       1987:599133 HCAPLUS Full-text
DOCUMENT NUMBER:
                       107:199133
ORIGINAL REFERENCE NO.: 107:31967a,31970a
TITLE:
                       Emulsifiers for emulsion polymerization
INVENTOR(S):
                       Oka, Masashi; Komiya, Kaoru
                      Asahi Denka Kogyo K. K., Japan
PATENT ASSIGNEE(S):
                       Jpn. Kokai Tokkyo Koho, 7 pp.
SOURCE:
                       CODEN: JKXXAF
DOCUMENT TYPE:
                       Patent
LANGUAGE:
                       Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:
    PATENT NO.
                  KIND
                                        APPLICATION NO.
                              DATE
                                          ______
                              19870515
    JP 62104802
                       Α
                                         JP 1986-171816
                                                                19860723
                                                 <--
    JP 05075001
                       В
                             19931019
                                         JP 1994-210055
    JP 07165806
                       Α
                              19950627
                                                                19940902
                                                <--
                       A
    JP 07165807
                             19950627
                                          JP 1994-210056
                                                                19940902
                                                <--
PRIORITY APPLN. INFO.:
                                          JP 1985-162055 A1 19850724
                                                 <--
ED
    Entered STN: 27 Nov 1987
AΒ
     The reactive title emulsifiers CH2:CRCH2OCH2CH(OX)CH2O(AO)nR1 (A = C2-4 alkylene;
     R = H, Me; R1 = C8-24 hydrocarbyl, acyl; n = 0-50; X = H, nonionic or nonionic
     hydrophilic group). Thus, nonylphenol was treated with allyl glycidyl ether in
     the presence of NaOH at 90.degree. for 5 h and ethoxylated to give an emulsifier
     (I). A soln. was prepd. from 80 g Et acrylate and 4 g I, and 8.4 g of this soln.
     was heated to 50.degree. with 0.08 g K2S2O8 and 0.04 g NaHSO4 in 120 g water to
     initiate the polymn., treated over 2 h with the remaining monomer soln. and further
     polymd. for 2 h to give a stable emulsion forming water-, weather- and
     heat-resistant films.
IT
    111115-38-9 111115-40-3
                              111123-58-1
                              111123-62-7
    111123-59-2 111123-61-6
    111123-64-9 111123-65-0 111144-58-2
    111144-59-3
```

(emulsifiers, reactive, for acrylic emulsion polymn.)

111115-38-9 HCAPLUS

RN

CN 2-Propanol, 1-(nonylphenoxy)-3-(2-propenyloxy)-, hydrogen sulfate, ammonium salt (9CI) (CA INDEX NAME)



$$D1 - (CH_2) 8 - Me$$

■ NH3

RN 111115-40-3 HCAPLUS

CN 2-Propanol, 1-(nonylphenoxy)-3-(2-propenyloxy)-, hydrogen sulfate, sodium salt (9CI) (CA INDEX NAME)



$$D1 - (CH2) 8 - Me$$

● Na

RN 111123-58-1 HCAPLUS

CN 2-Propanol, 1-(dodecyloxy)-3-(2-propenyloxy)-, dihydrogen phosphate, dipotassium salt (9CI) (CA INDEX NAME)

●2 K

10/596,747

RN 111123-59-2 HCAPLUS

CN 2-Propanol, 1-(2-propenyloxy)-3-(tridecyloxy)-, dihydrogen phosphate, dipotassium salt (9CI) (CA INDEX NAME)

●2 K

RN 111123-61-6 HCAPLUS

CN 2-Propanol, 1-(dodecyloxy)-3-(2-propenyloxy)-, dihydrogen phosphate, disodium salt (9CI) (CA INDEX NAME)

•2 Na

RN 111123-62-7 HCAPLUS

CN 2-Propanol, 1-(2-propenyloxy)-3-(tridecyloxy)-, dihydrogen phosphate, disodium salt (9CI) (CA INDEX NAME)

●2 Na

RN 111123-64-9 HCAPLUS

CN 2-Propanol, 1-(dodecyloxy)-3-(2-propen-1-yloxy)-, 2-(hydrogen sulfate), sodium salt (1:1) (CA INDEX NAME)

$$\begin{array}{c} \text{OSO3H} \\ \text{Me- (CH2)} \\ \text{11-O-CH2-CH-CH2-O-CH2-CH=CH2} \end{array}$$

Na

RN 111123-65-0 HCAPLUS

CN 2-Propanol, 1-(2-propen-1-yloxy)-3-(tridecyloxy)-, 2-(hydrogen sulfate), sodium salt (1:1) (CA INDEX NAME)

$$\begin{array}{c} \text{OSO3H} \\ \text{Me- (CH2)} \ 12 - \text{O-CH2} - \text{CH- CH2- O- CH2- CH- CH2-} \end{array}$$

● Na

RN 111144-58-2 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-sulfo-.omega.-[1[(nonylphenoxy)methyl]-2-(2-propen-1-yloxy)ethoxy]-, sodium salt (1:1)
(CA INDEX NAME)



 $D1 - (CH_2) 8 - Me$

Na

RN 111144-59-3 HCAPLUS

CN Poly(oxy-1,2-ethanediyl), .alpha.-[3-(2-propenyloxy)-2-(sulfooxy)propyl]-.omega.-(nonylphenoxy)-, sodium salt (9CI) (CA INDEX NAME)



 $D1 - (CH_2) 8 - Me$

Na

IT 111123-60-5 111123-63-8 111123-66-1

111123-67-2 111123-68-3 111123-69-4

111144-68-4 111144-69-5 111144-72-0

1111144-73-1 1111165-72-1 1111165-73-2

111165-74-3 111165-75-4

(emulsions, self-emulsifying, for water-, weather- and heat-resistant films)

RN 111123-60-5 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with 1-[(dodecyloxy)methyl]-2-(2-propenyloxy)ethyl dipotassium phosphate, ethenylbenzene and 1-[(2-propenyloxy)methyl]-2-(tridecyloxy)ethyl dipotassium phosphate (9CI) (CA INDEX NAME)

CM 1

CRN 111123-59-2 CMF C19 H39 O6 P . 2 K

OPO3H2 | Me- (CH2) 12-O-CH2-CH-CH2-O-CH2-CH=CH2

●2 K

CM 2

CRN 111123-58-1

CMF C18 H37 O6 P . 2 K

●2 K

CM 3

CRN 141-32-2 CMF C7 H12 O2

CM 4

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 111123-63-8 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with disodium
1-[(dodecyloxy)methyl]-2-(2-propenyloxy)ethyl phosphate, disodium
1-[(2-propenyloxy)methyl]-2-(tridecyloxy)ethyl phosphate and
ethenylbenzene (9CI) (CA INDEX NAME)

CM 1

CRN 111123-62-7

CMF C19 H39 O6 P . 2 Na

$$\begin{array}{c} \text{OPO3H2} \\ \text{Me- (CH2) } 12 - \text{O-CH2} - \text{CH- CH2- O- CH2- CH- CH2} \end{array}$$

•2 Na

CM 2

CRN 111123-61-6

CMF C18 H37 O6 P . 2 Na

●2 Na

CM 3

CRN 141-32-2 CMF C7 H12 O2

CM 4

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 111123-66-1 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with ethenylbenzene, sodium 1-[(dodecyloxy)methyl]-2-(2-propenyloxy)ethyl sulfate and sodium 1-[(2-propenyloxy)methyl]-2-(tridecyloxy)ethyl sulfate (9CI) (CA INDEX NAME)

CM 1

CRN 111123-65-0

CMF C19 H38 O6 S . Na

● Na

CM 2

CRN 111123-64-9 CMF C18 H36 O6 S . Na

QSO3H

● Na

CM 3

CRN 141-32-2 CMF C7 H12 O2

CM 4

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 111123-67-2 HCAPLUS

CN 2-Propenoic acid, ethyl ester, polymer with dipotassium
1-[(dodecyloxy)methyl]-2-(2-propenyloxy)ethyl phosphate and
dipotassium 1-[(2-propenyloxy)methyl]-2-(tridecyloxy)ethyl phosphate
(9CI) (CA INDEX NAME)

CM 1

CRN 111123-59-2

CMF C19 H39 O6 P . 2 K

 $\begin{array}{c} \text{OPO3H2} \\ \text{Me- (CH2)} \\ \text{12-O-CH2-CH-CH2-O-CH2-CH=-CH2} \end{array}$

●2 K

CM 2

CRN 111123-58-1 CMF C18 H37 O6 P . 2 K

OPO3H2 Me- (CH2) 11-O-CH2-CH-CH2-O-CH2-CH=CH2

●2 K

CM 3

CRN 140-88-5 CMF C5 H8 O2

0 || Eto-C-CH-CH2

RN 111123-68-3 HCAPLUS

CN 2-Propenoic acid, ethyl ester, polymer with disodium
1-[(dodecyloxy)methyl]-2-(2-propenyloxy)ethyl phosphate and disodium
1-[(2-propenyloxy)methyl]-2-(tridecyloxy)ethyl phosphate (9CI) (CA INDEX NAME)

CM 1

CRN 111123-62-7 CMF C19 H39 O6 P . 2 Na

$$\begin{array}{c} \text{OPO3H2} \\ \text{Me- (CH2)} \ 12 - \text{O-CH2} - \text{CH-CH2} - \text{O-CH2} - \text{CH-CH2} \\ \end{array}$$

•2 Na

CM 2

CRN 111123-61-6 CMF C18 H37 O6 P . 2 Na

$$\begin{array}{c} \text{OPO3H2} \\ \text{Me- (CH2)} \\ \text{11-O-CH2-CH-CH2-O-CH2-CH=-CH2} \end{array}$$

●2 Na

CM 3

CRN 140-88-5 CMF C5 H8 O2

RN 111123-69-4 HCAPLUS

CN 2-Propenoic acid, ethyl ester, polymer with sodium
1-[(dodecyloxy)methyl]-2-(2-propenyloxy)ethyl sulfate and sodium
1-[(2-propenyloxy)methyl]-2-(tridecyloxy)ethyl sulfate (9CI) (CA INDEX NAME)

CM 1

CRN 111123-65-0 CMF C19 H38 O6 S . Na

$$\begin{array}{c} \text{OSO3H} \\ \text{Me- (CH2)} \ 12 - \text{O-CH2} - \text{CH- CH2- O- CH2- CH== CH2} \end{array}$$

● Na

CM 2

CRN 111123-64-9 CMF C18 H36 O6 S . Na

$$\begin{array}{c} \text{OSO3H} \\ \text{Me- (CH2)} \\ \text{11-O-CH2-CH-CH2-O-CH2-CH=CH2} \end{array}$$

● Na

CM 3

CRN 140-88-5 CMF C5 H8 O2

RN 111144-68-4 HCAPLUS
CN 2-Propenoic acid, ethyl ester, polymer with
 .alpha.-[3-(2-propenyloxy)-2-(sulfooxy)propyl]-.omega. (nonylphenoxy)poly(oxy-1,2-ethanediyl) sodium salt (9CI) (CA INDEX NAME)

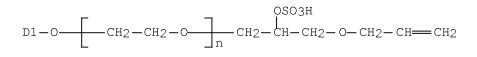
CM 1

CRN 111144-59-3

CMF (C2 H4 O)n C21 H34 O6 S . Na



 $D1 - (CH_2) 8 - Me$



● Na

CM 2

CRN 140-88-5 CMF C5 H8 O2

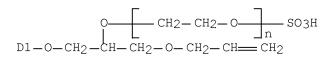
CM 1

CRN 111144-58-2

CMF (C2 H4 O)n C21 H34 O6 S . Na



$$D1 - (CH_2) 8 - Me$$



● Na

CM 2

CRN 140-88-5 CMF C5 H8 O2

RN 111144-72-0 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with ethenylbenzene and .alpha.-[3-(2-propenyloxy)-2-(sulfooxy)propyl]-.omega.(nonylphenoxy)poly(oxy-1,2-ethanediyl) sodium salt (9CI) (CA INDEX NAME)

CM 1

CRN 111144-59-3

CMF (C2 H4 O)n C21 H34 O6 S . Na $\,$



 $D1 - (CH_2) 8 - Me$

$${\tt D1-O-CH_2-CH_2-O-J_n-CH_2-CH-CH_2-O-CH_2-CH=CH_2}$$

● Na

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 111144-73-1 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with ethenylbenzene and .alpha.-sulfo-.omega.-[1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethoxy]poly(oxy-1,2-ethanediyl) sodium salt (9CI) (CA INDEX NAME)

CM 1

CRN 111144-58-2

CMF (C2 H4 O)n C21 H34 O6 S . Na



$$D1 - (CH_2) 8 - Me$$

• Na

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 100-42-5 CMF C8 H8

 H_2C \longrightarrow CH \longrightarrow Ph

RN 111165-72-1 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with ammonium 1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethyl sulfate and ethenylbenzene (9CI) (CA INDEX NAME)

CM 1

CRN 111115-38-9

CMF C21 H34 O6 S . H3 N

CCI IDS



$$D1 - (CH_2) 8 - Me$$

● инз

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 111165-73-2 HCAPLUS

CN 2-Propenoic acid, butyl ester, polymer with ethenylbenzene and sodium 1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethyl sulfate (9CI) (CA INDEX NAME)

CM 1

CRN 111115-40-3

CMF C21 H34 O6 S . Na

CCI IDS



$$D1 - (CH2) 8 - Me$$

● Na

CM 2

CRN 141-32-2 CMF C7 H12 O2

CM 3

CRN 100-42-5 CMF C8 H8

H2C=CH-Ph

RN 111165-74-3 HCAPLUS

CN 2-Propenoic acid, ethyl ester, polymer with ammonium 1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethyl sulfate (9CI) (CA INDEX NAME)

CM 1

CRN 111115-38-9

CMF C21 H34 O6 S . H3 N

CCI IDS



$$D1 - (CH2) 8 - Me$$

● NH3

CM 2

CRN 140-88-5 CMF C5 H8 O2

RN 111165-75-4 HCAPLUS

CN 2-Propenoic acid, ethyl ester, polymer with sodium
1-[(nonylphenoxy)methyl]-2-(2-propenyloxy)ethyl sulfate (9CI) (CA INDEX NAME)

CM 1

CRN 111115-40-3

CMF C21 H34 O6 S . Na

CCI IDS



$$D1 - (CH_2) 8 - Me$$

● Na

```
CM
         2
    CRN 140-88-5
    CMF C5 H8 O2
Eto-C-CH=CH2
IPCI C08F0002-24 [ICM,4]; B01F0017-42 [ICS,4]
IPCR B01F0017-42 [I,A]; B01F0017-00 [I,A]; C08F0002-24 [I,A]; C08F0216-14
    [N,A]; C08F0230-02 [N,A]
    35-4 (Chemistry of Synthetic High Polymers)
CC
    Section cross-reference(s): 46
IT
    25322-68-3D, ethers with secondary alcs., allyl derivs. 111100-57-3
    111100-58-4 111115-38-9
                              111115-40-3
    111123-58-1 111123-59-2 111123-61-6
    111123-62-7 111123-64-9
                              111123-65-0
    111144-58-2 111144-59-3
                              111144-60-6
    111144-61-7
        (emulsifiers, reactive, for acrylic emulsion polymn.)
    111100-59-5 111100-60-8
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                               111123-67-2
    111123-63-8
                 111123-66-1
    111123-68-3 111123-69-4
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    111144-67-3 111144-68-4 111144-69-5
    111144-70-8 111144-71-9 111144-72-0
    111144-73-1 111165-72-1 111165-73-2
                111165-75-4
    111165-74-3
        (emulsions, self-emulsifying, for water-, weather- and
       heat-resistant films)
OS.CITING REF COUNT:
                              THERE ARE 2 CAPLUS RECORDS THAT CITE THIS
                              RECORD (3 CITINGS)
```

=> d que 148

L5 STR

SO3H@6 G17 CH2



A @13

G3 17

VAR G1=1/6 VAR G2=O/N/S VAR G3=8/12 NODE ATTRIBUTES: DEFAULT MLEVEL IS ATOM

DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

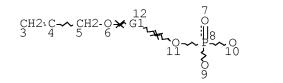
RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 17

STEREO ATTRIBUTES: NONE

L7 28963 SEA FILE=REGISTRY SSS FUL L5

L10 STR



REP G1=(1-20) 13 NODE ATTRIBUTES:

NSPEC IS RC AT 13
DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 11

STEREO ATTRIBUTES: NONE

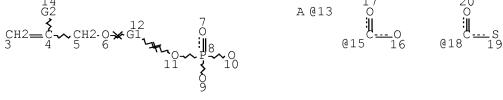
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L16 STR

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GRAPH ATTRIBUTES: RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 13

STEREO ATTRIBUTES: NONE 13442 SEA FILE=REGISTRY SUB=L7 SSS FUL L16 L18 14 SEA FILE=REGISTRY SPE=ON ABB=ON PLU=ON L12 AND L18 L21 8 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L20 L22 400 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L12 13991 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L18 L23 L25 58 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L22 AND PHARM?/SC. SX L26 44 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L25 AND (1802-2003)/PRY,AY,PY L27 37 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L22 AND L23 L28 20 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L27 AND (1802-2003)/PRY,AY,PY 2 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L21 AND (1802-2003 L29)/PRY,AY,PY L30 20 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L28 OR L29 L33 STR



23 0 11 0 21 22

REP G1=(1-20) 13 VAR G2=15/18/21/COOH/SO3H NODE ATTRIBUTES: NSPEC IS RC AT 13 DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 21

STEREO ATTRIBUTES: NONE

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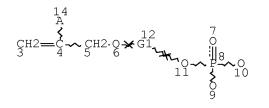
L37 8 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L35

L38 3 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L37 AND (1802-2003

)/PRY,AY,PY

L39 19 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L30 NOT L38

L40 STF



A@13

REP G1=(1-20) 13

NODE ATTRIBUTES:

NSPEC IS RC AT 13 DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 12

STEREO ATTRIBUTES: NONE

L42

112 SEA FILE=REGISTRY SUB=L12 SSS FUL L40

L43

30 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L42

L45

15 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L43 AND (1802-2003)

)/PRY,AY,PY

L46

54 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L45 OR L26

L47

48 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L46 NOT (38 OR L39)

L48

39 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L47 AND PHARM3/SC

L48 39 SEA FILE=HCAPLUS SPE=ON ABB=ON PLU=ON L47 AND PHARM?/SC,

\Rightarrow d 148 1-39 ibib ed abs fhitstr hitind

L48 ANSWER 1 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2010:1405036 HCAPLUS Full-text

DOCUMENT NUMBER: 153:618952

TITLE: Novel hexahydropyrazinotriazinedione compounds of

reverse turn mimetics and their preparation and

use thereof

INVENTOR(S): Chung, Jae Uk; Jung, Kyung-Yun; Jeong, Min-Wook;

Jung, Hee-Kyung; La, Hyun-Ju

PATENT ASSIGNEE(S): Choongwae Pharma Corporation, S. Korea

U.S. Pat. Appl. Publ., 320 pp., Cont.-in-part of U.S. Ser. No. 974,941. SOURCE:

CODEN: USXXCO

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 11

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US	2004	0072	831		A1		2004	0415		US 2	003-	 4118 	77		20	0030409
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	7232 2007		425		B2 A1		2007 2007			US 2	004-		72		20	0040416
	7576 2007		052		B2 A1		2009 2007			US 2	005-	1081	64		20	0050415
	7566 7671				B2 B1		2009 2010			US 2	007-		41		20	0071015
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SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR,

NE, SN, TD, TG, BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AP, EA, EP, OA PRIORITY APPLN. INFO.: US 2001-976470 B2 20011012 <--US 2002-87443 B2 20020301 <--WO 2002-KR1901 A 20021011 <--US 2003-411877 A2 20030409 <--US 2004-803179 A2 20040317

> US 2004-826972 A2 20040416 US 2005-108164 A2 20050415 US 2007-974941 A2 20071015

WO 2008-KR6070 20081015 CN 2002-822567 A3 20021011

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ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 153:618952 Entered STN: 12 Nov 2010

GΙ

AΒ Conformationally constrained compds. of formula I that mimic the secondary structure of reverse-turn regions of biol. active peptides and proteins and having bicyclic framework are disclosed, as well as their prodrugs. Compds. of formula I wherein E is CO, CHR3 and NR3; W is CO, CONH, CO2, COS, SO2 and a bond; R1 is (un) substituted (hetero) aryl; R2, R3, R4 and R5 are independently an amino acid side chain moiety or an amino acid side chain deriv. and stereoisomers, mixt. of stereoisomers and pharmaceutically acceptable salts thereof, are claimed. Such reverse-turn mimetic structures and prodrugs have utility over a wide range of fields, including use as diagnostic and therapeutic agents. The invention also relates to a use of such compds. for the prepn. of a medicament for treating or

preventing cancer including an acute myeloid leukemia. Example compd. II was prepd. by a general procedure (general procedure given). All the invention compds. were evaluated for their CYP3A4 inhibitory activity (some data given).

IT 1145676-59-0P

(drug candidate; prepn. of hexahydropyrazinotriazinedione compds. as CYP3A4 inhibitors useful as reverse turn mimetics and in the treatment of cancer)

RN 1145676-59-0 HCAPLUS

CN 2H-Pyrazino[2,1-c][1,2,4]triazine-1(6H)-carboxamide, hexahydro-4,7-dioxo-N-(phenylmethyl)-6-[[4-(phosphonooxy)phenyl]methyl]-2-(2-propen-1-yl)-8-[[4-(2-propen-1-yloxy)phenyl]methyl]-, sodium salt (1:2), (6S,9aS)- (CA INDEX NAME)

Absolute stereochemistry.

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INCL 514081000; 544184000; 544112000; 514243000; 514233200
IPCI A61K0031-53 [I,A]; C07D0487-04 [I,A]; A61K0031-5377 [I,A];
     A61K0031-675 [I,A]; A61P0035-02 [I,A]; C07D0487-04 [I,A]; A61K0031-53
     [I,A]; A61P0035-00 [I,A]
     514/081.000; 514/233.200; 514/243.000; 544/112.000; 544/184.000
NCL
     28-19 (Heterocyclic Compounds (More Than One Hetero Atom))
CC
     Section cross-reference(s): 1, 34, 63
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(drug candidate; prepn. of hexahydropyrazinotriazinedione compds. as CYP3A4 inhibitors useful as reverse turn mimetics and in the treatment of cancer)

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L48 ANSWER 2 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2010:261208 HCAPLUS Full-text DOCUMENT NUMBER: 152:335464
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TITLE: Reverse-turn mimetics for treatment of cancer and

rheumatoid arthritis

INVENTOR(S): Moon, Sung Hwan; Chung, Jae Uk; Lee, Sung Chan; Equchi, Masakatsu; Kahn, Michael; Jeong, Kwang

Won; Nguyen, Cu; Lee, Soo Jin

PATENT ASSIGNEE(S): Choongwae Pharma Corp., S. Korea

SOURCE: U.S., 1320 pp., Cont.-in-part of U.S. Ser. No.

108,164.

CODEN: USXXAM

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 11

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ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT OTHER SOURCE(S): CASREACT 152:335464; MARPAT 152:335464 ED Entered STN: 03 Mar 2010 GI

The invention discloses conformationally-constrained compds. I [A is CO; B is CHR4; D is CO; E is Z-R6; G is X-R7; W is CONH, COO, COS, SO2 or null; Z = CH; X = N; R1-2, R4, R6-7 are the same or different and independently selected from an amino acid side chain or deriv.], their stereoisomers and pharmaceutically acceptable

II

10/596.747

salts that mimic the secondary structure of reverse-turn regions of biol. active peptides and proteins and have utility over a wide range of fields, including use as diagnostic and therapeutic agents. Libraries contg. the reverse-turn mimetic structures of this invention are also disclosed as well as methods for screening them to identify biol. active members. The invention also relates to the use of such compds. for inhibiting or treating disorders modulated by the Wnt-signaling pathway, such as cancer, restenosis assocd. with angioplasty, polycystic kidney disease, aberrant angiogenesis disease, rheumatoid arthritis, tuberous sclerosis complex, Alzheimer's disease, excess hair growth or loss, or ulcerative colitis. Thus, triazolopyrazinone deriv. II was prepd. using a bromoacetal resin and showed IC50 = 2.349 .mu.M against SW480 cells.

IT 1145676-59-0P

(prepn. of reverse-turn mimetics for treatment of cancer and rheumatoid arthritis)

RN 1145676-59-0 HCAPLUS

CN 2H-Pyrazino[2,1-c][1,2,4]triazine-1(6H)-carboxamide, hexahydro-4,7-dioxo-N-(phenylmethyl)-6-[[4-(phosphonooxy)phenyl]methyl]-2-(2-propen-1-yl)-8-[[4-(2-propen-1-yloxy)phenyl]methyl]-, sodium salt (1:2), (6S,9aS)- (CA INDEX NAME)

Absolute stereochemistry.

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INCL 514243000; 544184000
IPCI C07D0487-04 [I,A]; A61K0031-53 [I,A]; A61P0019-02 [I,A]; A61P0035-00
IPCR C07D0487-04 [I,A]; A61K0031-53 [I,A]; A61P0019-02 [I,A]; A61P0035-00
     [N,A]
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(prepn. of reverse-turn mimetics for treatment of cancer and rheumatoid arthritis)

REFERENCE COUNT: 149

THERE ARE 149 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L48 ANSWER 3 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2009:490612 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 150:447990

TITLE: Novel hexahydropyrazinotriazinedione compounds of

reverse turn mimetics and their preparation and

use thereof

INVENTOR(S): Chung, Jae Uk; Jung, Kyung-Yun; Jeong, Min-Wook;

Jung, Hee-Kyung; La, Hyun-Ju

PATENT ASSIGNEE(S): Choongwae Pharma Corporation, S. Korea

SOURCE: PCT Int. Appl., 192 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 11

PATENT NO.	KIND DATE	APPLICAT:	APPLICATION NO.					
WO 2009051397	A2 2009	00423 WO 2008-I	KR6070	20081015				
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ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT OTHER SOURCE(S): CASREACT 150:447990; MARPAT 150:447990 ED Entered STN: 24 Apr 2009

GΙ

Conformationally constrained compds. of formula I that are novel and mimic the secondary structure of reverse-turn regions of biol. active peptides and proteins and having bicyclic framework are disclosed, as well as their prodrugs.1 Compds. of formula I wherein E is CO, CHR3 and NR3; W is CO, CONH, CO2, COS, SO2 and a bond; R1 is (un) substituted (hetero) aryl; R2, R3, R4 and R5 are independently an amino acid side chain moiety or an amino acid side chain deriv. and stereoisomers, mixt. of stereoisomers and pharmaceutically acceptable salts thereof, are claimed. Such reverse-turn mimetic structures and prodrugs have utility over a wide range of fields, including use as diagnostic and therapeutic agents. The invention also relates to a use of such compds. for the prepn. of a medicament for treating or preventing cancer including an acute myeloid leukemia. Example compd. II was prepd. by a general procedure (general procedure given). All the invention compds. were evaluated for their CYP3A4 inhibitory activity (some data given).

IT 1145676-59-0P

(drug candidate; prepn. of hexahydropyrazinotriazinedione compds. as CYP3A4 inhibitors useful as reverse turn mimetics and in the treatment of cancer)

RN 1145676-59-0 HCAPLUS

CN 2H-Pyrazino[2,1-c][1,2,4]triazine-1(6H)-carboxamide, hexahydro-4,7-dioxo-N-(phenylmethyl)-6-[[4-(phosphonooxy)phenyl]methyl]-2-(2-propen-1-yl)-8-[[4-(2-propen-1-yloxy)phenyl]methyl]-, sodium salt (1:2), (6S,9aS)- (CA INDEX NAME)

Absolute stereochemistry.

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IPCI C07D0487-04 [I,A]; C07D0487-00 [I,C]; C07D0487-04 [I,A]; C07D0487-00
     [I,C]; C07D0487-04 [I,A]
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     28-19 (Heterocyclic Compounds (More Than One Hetero Atom))
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       (drug candidate; prepn. of hexahydropyrazinotriazinedione compds.
       as CYP3A4 inhibitors useful as reverse turn mimetics and in the
       treatment of cancer)
OS.CITING REF COUNT:
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L48 ANSWER 4 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN
ACCESSION NUMBER:
                       2005:568976 HCAPLUS Full-text
                       143:83603
                       One-part self-etching, self-priming dental
                       adhesive composition
                       Klee, Joachim E.; Lehmann, Uwe; Walz, Uwe
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DOCUMENT NUMBER:

TITLE:

INVENTOR(S):

PATENT ASSIGNEE(S): Dentsply Detrey GmbH, Germany

SOURCE: Eur. Pat. Appl., 30 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

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PRIORITY APPLN. INFO.:
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ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

ED Entered STN: 01 Jul 2005

One-part self-etching, self-priming dental adhesive compn. having a pH of at most 2 comprises (a) a polymerizable acidic phosphoric acid ester monomer; (b) one or more polymerizable acidic monomers; (c) a polymerizable N-substituted alkylacrylic or acrylic acid amide monomer; (d) an org. and/or inorg. acid; (e) an org. water sol. solvent and/or water; and (f) polymn. initiator, inhibitor and stabilizer. An adhesive polymer was prepd. from 2-acrylamido-2-methyl-propane-sulfonic acid, 3, (4),8, (9)-bis(acrylamido methyl) tricyclo-5.2.1.02,6 decane, Et 2-[13-dihydrogen phosphoryl-13,2-dioxatridecyl]acrylate, and N,N'-bisacrylamido-N,N'-diethyl-1,3-propane.

IT 752234-98-3P

(one-part self-etching, self-priming dental adhesive compn.)

RN 752234-98-3 HCAPLUS

CN 2-Propenoic acid, 2-[[[10-(phosphonooxy)decyl]oxy]methyl]-, 1-ethyl ester (CA INDEX NAME)

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(one-part self-etching, self-priming dental adhesive compn.)
    855894-57-4P, 2-Acrylamido-2-methyl-propane-sulfonic
ΙT
    acid-3, (4), 8, (9) -bis(acrylamido methyl) tricyclo-5.2.1.02,6
    decane-Ethyl 2-[13-dihydrogen phosphoryl-13,2-dioxatridecyl]acrylate-
    N, N'-Bisacrylamido-N, N'-diethyl-1, 3-propane copolymer
    855894-58-5P
        (one-part self-etching, self-priming dental adhesive compn.)
OS.CITING REF COUNT:
                              THERE ARE 7 CAPLUS RECORDS THAT CITE THIS
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REFERENCE COUNT:
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L48 ANSWER 5 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN
                       2005:216597 HCAPLUS Full-text
ACCESSION NUMBER:
                        142:291323
DOCUMENT NUMBER:
TITLE:
                        Compositions and methods for the treatment of
                        severe acute respiratory syndrome (SARS)
INVENTOR(S):
                       Hardee, Greg; Dellamary, Luis
                       Isis Pharmaceuticals, Inc., USA
PATENT ASSIGNEE(S):
SOURCE:
                        PCT Int. Appl., 217 pp.
                        CODEN: PIXXD2
DOCUMENT TYPE:
                        Patent
LANGUAGE:
                        English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:
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    WO 2005020885 A2
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PRIORITY APPLN. INFO.:
                                           US 2003-472774P P 20030521
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ED
    Entered STN: 11 Mar 2005
AΒ
     The invention provides compns. and methods for treating a coronavirus infection,
     esp. a SARS CoV infection. The compns. comprise an antiviral nucleoside or mimetic
     thereof, or an antiviral antisense agent, in a form suitable for pulmonary or nasal
     delivery. The methods comprise administration to a patient in need thereof the
     effective amt. of an antiviral compn. by pulmonary or nasal instillation.
IT
    847648-34-4
        (compns. and methods for treatment of severe acute respiratory
       syndrome)
    847648-34-4 HCAPLUS
RN
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CN Guanosine 5'-(tetrahydrogen triphosphate), 2'-O-2-propenyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

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IPCI A61K [ICM, 7]
IPCR A61K [I,S]; A61K0031-7052 [I,A]; C07H0019-22 [I,A]
CC
     1-5 (Pharmacology)
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ACCESSION NUMBER: 2005:120948 HCAPLUS Full-text
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DOCUMENT NUMBER:
TITLE:
                                                Preparation of nucleobase phosphonate analogs for
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                                                Krawczyk, Steven H.
INVENTOR(S):
                                             Gilead Sciences, Inc., USA
PATENT ASSIGNEE(S):
SOURCE:
                                                  PCT Int. Appl., 140 pp.
                                                  CODEN: PIXXD2
DOCUMENT TYPE:
                                                 Patent
LANGUAGE:
                                                  English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:
          PATENT NO.
                                                                                     APPLICATION NO.
                                               KIND DATE
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                                                                                        _____
                                                A2
                                                                20050210 WO 2004-US24922 20040730
         WO 2005012324
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          WO 2005012324
                                                 A3 20050506
                  W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA,
                          CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI,
                          GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP,
                          KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW,
                          MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD,
                          SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ,
                          VC, VN, YU, ZA, ZM, ZW
                  RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW,
                          AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ,
                          DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL,
                          PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ,
                          GW, ML, MR, NE, SN, TD, TG
                                                                                    AU 2004-260789
         AU 2004260789
                                                 A1 20050210
                                                                                                                                     20040730
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         AU 2004260789 B2 20110630
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CA	2533	966			A1	2005	0210	CA	2004-2		2	20040730		
US	2005	0059	637		A1	2005	0317	US	2004-9	90328	88		2	20040730
	7579 1656				B2 A2	2009 2006		EP	2004-7	77985	55		2	20040730
	R:	•	•	•	•	DK, ES, RO, CY,	•	•	•	LI,	•	•	SE,	MC,
JP	2007	•	•	~-,	T	2007	•	•	2006-5	52212	•		2	20040730
NZ	5449	88			A	2009	1127	NΖ	2004-5	54498	8 8		2	20040730
US	2006	0252	729		A1	2006	1109	US	2006-5	56681	. 9		2	20060127
PRIORIT	Y APP	LN.	INFO	.:				US	2003-4	49112	23P	1	2	20030730
								WO	2004-0		22	Ţ	N 2	20040730

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT OTHER SOURCE(S): CASREACT 142:197761; MARPAT 142:197761 ED Entered STN: 11 Feb 2005

GΙ

AΒ The present invention provides the prepn. of nucleobase phosphonate analogs I, wherein X1 is C:CR1R2, C:N-OR, C:O, CR1R2, C:N-NR2, S:O; X2 is O, NR, S; Y1 is O, S, NR, N-Nr2, +N(O)(R), +N(OR), +N(O)(OR), sulfonyl; Y3 and Z are independently H, OH, OR, NR2, CN, NO2, F, Cl, Br, iodo; Rx is H, protecting group; R1-R8 are independently H, F, Cl, Br, iodo, OH, C(:Y1)R, C(:Y1)OR, C(:Y1)NR2+NR3, SR, S(O)R, S(O2)R, S(O) 2(ORx), OC(:Y)Rx, -OC(=Y1)OR, -OC(=Y1)(N(Rx) 2), -SC(=Y1)Rx, -OC(=Y1)RxSC(=Y1)OR, -SC(=Y1)(N(Rx)2), -N(Rx)C(=Y1)R, -N(Rx)C(=Y1)OR, or -N(Rx)C(=Y1)ORN(Rx)C(=Y1)N(Rx)2, amino, ammonium, alkylamino, dialkylamino, trialkyl-ammonium, alkyl, alkyl-halide, carboxylate sulfate, sulfamate, sulfonate, 5-7 membered ring sultam, alkyl sulfonate, alkylamino, 4-dialkylamino-pyridinium, alkyl-hydroxy, alkyl-thiol, alkyl sulfone, aryl sulfone, aryl sulfoxide, arylthio, sulfonamide, alkyl sulfoxide, ester, amido, 5-7 membered ring lactam, 5-7 membered ring lactone, nitrile, azido, nitro, alkoxy, alkyl, alkenyl, alkynyl, aryl, heteroaryl, polyethylene-oxy; two of R1-R8 form a carbocyclic ring of 3 to 7 carbon atoms; R is alkyl, alkenyl, alkynyl, aryl, with activity against infectious viruses. The

Ι

compds. of the invention may inhibit retroviral reverse transcriptases and thus inhibit the replication of the virus. They are useful for treating human patients infected with a human retrovirus, such as human immunodeficiency virus (strains of HIV-1 or HIV-2) or human T-cell leukemia viruses (HTLV-I or HTLV-II) which results in acquired immunodeficiency syndrome (AIDS) and/or related diseases. The present invention also relates generally to the accumulation or retention of therapeutic compds. inside cells. The invention is more particularly related to attaining high concns. of active metabolite mols. in HIV infected cells. Intracellular targeting may be achieved by methods and compns. which allow accumulation or retention of biol. active agents inside cells. Such effective targeting may be applicable to a variety of therapeutic formulations and procedures. Thus, nucleobase phosphonate II was prepd. and tested as antiviral agent and retroviral reverse transcriptase inhibitor. Within the context of the invention, typically compns. are first screened for inhibition of HIV reverse transcriptase in vitro and compns. showing inhibitory activity are then screened for activity in vivo. Compns. having in vitro Ki (inhibitory consts.) of less then about 5 \times 10-6 M, typically less than about 1 \times 10-7 M and preferably less than about 5 \times 10-8 \times are preferred for in vivo use.

IT 839711-04-5P

(prepn. of nucleobase phosphonate analogs for antiviral treatment and as retroviral reverse transcriptase inhibitors)

RN 839711-04-5 HCAPLUS

CN Diphosphoric acid, monoanhydride with

[[3-(3-amino-1,6-dihydro-6-oxo-9H-purin-9-yl)-4-hydroxy-2-methylenebutoxy]methyl]phosphonic acid (9CI) (CA INDEX NAME)

$$H_{2N}$$
 H_{2N}
 H

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IPCI C07H0019-00 [ICM,7]
IPCR C07F0009-6561 [I,A]
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CC 26-9 (Biomolecules and Their Synthetic Analogs)

Section cross-reference(s): 1, 7, 33, 63

839710-27-9P 839710-29-1P 839710-43-9P 839710-45-1P ΙT 839710-56-4P 839710-59-7P 839710-73-5P 839710-77-9P 839710-83-7P 839710-86-0P 839710-78-0P 839710-85-9P 839710-90-6P 839710-93-9P 839711-02-3P 839711-04-5P 839711-13-6P 839711-18-1P

839711-08-9P 839711 839711-19-2P

(prepn. of nucleobase phosphonate analogs for antiviral treatment and as retroviral reverse transcriptase inhibitors)

OS.CITING REF COUNT: 2 THERE ARE 2 CAPLUS RECORDS THAT CITE THIS RECORD (2 CITINGS)

REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L48 ANSWER 7 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN

ACCESSION NUMBER: 2004:732258 HCAPLUS Full-text

DOCUMENT NUMBER: 141:243056

TITLE: Polymerizable phosphoric acid ester derivatives

for dental compositions

INVENTOR(S): Klee, Joachim E.; Lehmann, Uwe; Walz, Uwe; Liu,

Huaibing

PATENT ASSIGNEE(S): Dentsply Detrey GmbH, Germany

SOURCE: Eur. Pat. Appl., 20 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

	PATENT NO.						KIND		DATE		APPLICATION NO.						DATE	
	EP	1454				A1	_	2004	0908	-	EP 2	 003- >				20030307		
		R:	•	•	•	•	•	ES, FI,	•	•	•	•	•	•	•	•	MC, HU, SK	
	CA	2518	202			A1		2004	0916	(CA 2		2518 	202		2	0040305	
	WO	2004	0781	00		A2		2004	0916	WO 2004-EP2289 <							0040305	
	WO	2004	0781	00		A3		2004	1028									
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			KR,		LC,	LK,		LS,										
		RW:	BE,	BG,	CH,	CY,	CZ,	MW, DE, PT,	DK,	EE,	ES,	FI,	FR,	GB,	GR,	HU,	IE,	
								GW,							ъо,	Cr,	cg,	
	ΕP	1601		011,	011,	A2		2005				004-				2	0040305	
	ΕP	1601	679			В1		2011	0511									
		R:	AT,	IE,				ES, FI,										
	JP	2006	5203	4 4		Т		2006	0907	!	JP 2		5045 	63		2	0040305	
	JΡ	4594	297			В2		2010	1208									
	US 20060246017				A1		2006	1102							2	0060626		
PRIOF	RIT	APP:	LN.	INFO	.:					:	EP 2		5174 		j	A 2	0030307	
										1	WO 2	004-	EP22	89	1	W 2	0040305	

ED Entered STN: 09 Sep 2004

AB The present invention provides a polymerizable phosphoric acid ester deriv. for use in dental compns. E.g.,

^{2,2,2}-tris(2,6-dioxa-4-methylene-5-oxo-octyl)ethanol phosphoric acid ester was prepd. from pentaerythritol, Et chloromethyacrylate, and then treatment with the product with POC13 and hydrolyzed.

IT 752234-96-1P

(polymerizable phosphoric acid ester derivs. for dental compns.) RN $\,$ 752234-96-1 HCAPLUS

CN 2-Propenoic acid, 2,2'-[[2-[[2-(ethoxycarbonyl)-2-propenyl]oxy]methyl]-2-[(phosphonooxy)methyl]-1,3-propanediyl]bis(oxymethylene)]bis-, 1,1'-diethyl ester (9CI) (CA INDEX NAME)

IPCI C07F0009-09 [ICM,7]; A61K0006-08 [ICS,7]; C08F0030-02 [ICS,7]
IPCR A61K0006-00 [I,A]; A61K0006-08 [I,A]; A61K0006-083 [I,A]; C07F0009-09

[I,A]; C08F0030-02 [I,A]
23-17 (Aliphatic Compounds)

CC 23-17 (Aliphatic Compounds)
Section cross-reference(s): 63

IT 752234-96-19 752234-98-39 752235-00-0P

(polymerizable phosphoric acid ester derivs. for dental compns.)
OS.CITING REF COUNT: 3 THERE ARE 3 CAPLUS RECORDS THAT CITE THIS

RECORD (3 CITINGS)

REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 8 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2004:252299 HCAPLUS Full-text

DOCUMENT NUMBER: 140:283384

TITLE: Modulators of GTPases and modulator-resistant

enzymes and their uses in drug design and target

validation

INVENTOR(S): Shah, Kavita; Vincent, Fabien; Cuento, Maria A. PATENT ASSIGNEE(S): Irm, Llc, UK; Novartis Pharmaceuticals Corporation

SOURCE: PCT Int. Appl., 144 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PAT	rent	NO.			KIN	D	DATE	;	APPLICATION NO.						DATE		
WO	2004	0240	82		A2	_	2004	0325	1	WO 2	 1-003 !>	US28	594		21	0030910	
WO	2004	0240	82		АЗ		2009	0618									
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		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	ES,	FI,	GB,	GD,	
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		LC,	LK,	LR,	LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	
		NΙ,	NO,	NΖ,	OM,	PG,	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,	SG,	SK,	
		SL,	SY,	ТJ,	TM,	TN,	TR,	TT,	TZ,	UA,	UG,	US,	UΖ,	VC,	VN,	YU,	

ZA, ZM, ZW RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG, AP, EA, EP, OA AU 2003267145 20040430 AU 2003-267145 Α1 20030910 <--US 2003-660113 US 20040241706 Α1 20041202 20030910 <--PRIORITY APPLN. INFO.: US 2002-410536P 20020913 Р <--US 2003-461755P 20030409 <--WO 2003-US28594 20030910 W <--

OTHER SOURCE(S): CASREACT 140:283384; MARPAT 140:283384

ED Entered STN: 26 Mar 2004

GΙ

AB Guanine derivs. that act as modulators of GTPases and GTPase variants that do not interact with these modulators are described for use in the design of improved modulators of GTPase activity. The method involves generating variants of the enzyme that do not interact with a known modulator and then developing effectors that interact with the resistant variant. The prepn. of guanosine derivs. and of a series of p21c-Ha-ras protein substitution variants is described,.

IT 674796-08-8P

(prepn. and use of; modulators of GTPases and modulator-resistant enzymes and their uses in drug design and target validation)

RN 674796-08-8 HCAPLUS

CN Guanosine 5'-(tetrahydrogen triphosphate), 6-0-2-propenyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IPCI A01N0043-02 [I,C]; A01N0043-04 [I,A]; A61K0031-70 [I,C]; A61K0031-70
 [I,A]
IPCR A01N0043-04 [I,A]; A61K0031-70 [I,A]; A61K0048-00 [I,A]; C07H0021-04
 [I,A]; C12Q0001-68 [I,A]

CC 7-5 (Enzymes)

Section cross-reference(s): 1, 3

1074-41-5DP, derivs. 15867-02-4P 17670-19-8P 26775-35-9P ΙT 26783-32-4P 26783-36-8P 67831-83-8DP, derivs. 99404-63-4P 674795-56-3P 282531-50-4P 674795-54-1P 674795-55-2P 674795-57-4P 674795-58-5P 674795-59-6P 674795-60-9P 674795-61-0P 674795-62-1P 674795-63-2P 674795-64-3P 674795-65-4P 674795-66-5P 674795-67-6P 674795-68-7P 674795-69-8P 674795-70-1P 674795-71-2P 674795-72-3P 674795-73-4P 674795-74-5P 674795-75-6P 674795-77-8P 674795-78-9P 674795-85-8P 674795-86-9P 674795-87-0P 674795-89-2P 674795-90-5P 674795-91-6P 674795-88-1P 674795-92-7P 674795-93-8P 674795-95-0P 674795-96-1P 674795-97-2P 674795-98-3P 674795-99-4P 674796-00-0P 674796-01-1P 674796-02-2P 674796-03-3P 674796-04-4P 674796-05-5P 674796-06-6P 674796-07-7P 674796-08-8P 674796-09-9P 674796-10-2P 674796-11-3P 674796-12-4P 674796-13-5P 674796-14-6P 674796-16-8P

(prepn. and use of; modulators of GTPases and modulator-resistant enzymes and their uses in drug design and target validation)

L48 ANSWER 9 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2003:972032 HCAPLUS Full-text

DOCUMENT NUMBER: 140:16928

TITLE: Synthetic methods for the large scale production

from glucose of analogs of sphingosine,

azidosphingosine, ceramides, lactosyl ceramides,

and glycosyl phytosphingosine

INVENTOR(S): Bundle, David R.; Ling, Chang Chun; Zhang, Ping

PATENT ASSIGNEE(S): Can.

SOURCE: PCT Int. Appl., 69 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003101937	A1	20031211	WO 2003-CA832	20030602

<--W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG AU 2003233725 Α1 20031219 AU 2003-233725 20030602 <--PRIORITY APPLN. INFO.: US 2002-384435P 20020531 <--WO 2003-CA832 20030602 <--CASREACT 140:16928; MARPAT 140:16928 OTHER SOURCE(S):

ED Entered STN: 14 Dec 2003

GΙ

AB Simple, direct and easily scaled synthetic methods were disclosed for the prodn. of sphingosines, azidosphingosines, ceramides, lactosyl ceramides, glycosyl phytosphingosines, such as ROCH2CH(R2)CH(OH)CH2CCH2CH:CH2 and ROCH2CH(R2)CH(OH)CH:CH2 [R = H, PO3H2, alkyl, alkenyl, alkynyl, aryl, heteroaryl, heterocyclyl, acyl, thioacyl, peptidyl, glycosyl; R2 = N2, NH2, acylamino, etc.], for use in pharmaceutical applications (no biol. testing data presented). The prepd. sphingosine and ceramide derivs. are useful intermediates for prepg. protein conjugates that could cause an immune response, and thus, may be useful for prepg. anti-cancer vaccines. Thus, (2S,3R)-2-azidopent-4-ene-1,3-diol was prepd. via a synthetic sequence which started from 1,2-0-isopropylidene-.alpha.-D-glucofuranose. Ultimately, the mono-acetate salt

of glycylated sphingosine analog I was prepd. I could be used to further link with a protein form a protein conjugate suitable as vaccine.

Ι

IT 631090~29~4P

(claimed compd.; synthetic methods for the large scale prodn. from glucose of analogs of sphingosine, azidosphingosine, ceramides, lactosyl ceramides, and glycosyl phytosphingosines)

RN 631090-29-4 HCAPLUS

CN 1,3-Butanediol, 2-azido-4-(2-propen-1-yloxy)-, 1-(dihydrogen phosphate), (2S,3S)- (CA INDEX NAME)

Absolute stereochemistry.

IPCI C07C0247-08 [ICM,7]; C07C0215-24 [ICS,7]; C07C0233-18 [ICS,7];

C07C0323-25 [ICS,7]

IPCR C07C0215-24 [I,A]; C07C0233-18 [I,A]; C07C0247-04 [I,A]; C07C0247-08
[I,A]; C07C0323-25 [I,A]; C07H0015-10 [I,A]

CC 33-7 (Carbohydrates)

Section cross-reference(s): 1, 15, 34, 63

IT 631090-29-4P 631090-30-7P 631090-31-8P 631090-32-9P

631090-33-0P

(claimed compd.; synthetic methods for the large scale prodn. from glucose of analogs of sphingosine, azidosphingosine, ceramides, lactosyl ceramides, and glycosyl phytosphingosines)

OS.CITING REF COUNT: 10 THERE ARE 10 CAPLUS RECORDS THAT CITE THIS

RECORD (12 CITINGS)

REFERENCE COUNT: 8 THERE ARE 8 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 10 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2003:656778 HCAPLUS Full-text

DOCUMENT NUMBER: 139:180298

TITLE: Preparation of substituted inositols and their use

as phosphatidylinositol hexamannoside mimics and

potential drug delivery agents

INVENTOR(S): Rademacher, Thomas William; Schmidt, Richard;

Stadelmaier, Andreas

PATENT ASSIGNEE(S): Lascaux Pharmaceuticals Limited, UK

SOURCE: PCT Int. Appl., 87 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT NO	Ο.	K	KIND	DATE			APPL:	ICAT:		DATE			
WO 200306	68789	-	A1	20030	Ţ	WO 20)-800 ->		20030213				
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(CN, CO,	CR, C	CU, CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	ES,	FΙ,	GB,	GD,
(GE, GH,	GM, H	HR, HU,	ID,	IL,	IN,	IS,	JP,	KΕ,	KG,	KΡ,	KR,	KΖ,
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E	BY, KG,	KZ, M	ΔD, RU,	ТJ,	TM,	AT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,

EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG 20030904 AU 2003-245767 AU 2003245767 Α1 20030213 <--EP 2003-739562 EP 1480991 Α1 20041201 20030213 <--R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK US 20050143290 Α1 20050630 US 2005-504605 <--GB 2002-3535 PRIORITY APPLN. INFO.: A 20020214 <--WO 2003-GB604 W 20030213 <--

OTHER SOURCE(S): MARPAT 139:180298

ED Entered STN: 22 Aug 2003

GΙ

HO.
$$R^4$$
 R^3
 R^6
 R^2
 R^3
 R^4
 R^3
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 R^3
 R^4
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 R^4
 R^3
 R^4
 R^3

AΒ Inositol phosphate esters and conjugates I and II, wherein R1 is hydroxyl, phosphate, phosphatidic acid or a phosphate ester; R2 is a sugar moiety; R3 is are selected from hydroxyl or phosphate; R4 and/or R6 is or are independently selected from: an amino acid; or a peptide or polypeptide; or a group having the general formula: O-(CH2)n-CH(NR7R8)-CO2X, wherein: n is an integer between 1 and 10, R7 and R8 are independently selected from hydrogen, nitrogen, acyl or alkyl; and X is hydrogen, alkyl or a cation where the terminal group is CO2-; or a substituted or unsubstituted arom. group, formed between the compds. and a coupling partner are disclosed, in particular compds. based on a myo-inositol which is substituted at position 1 with a phosphate ester group, at position 2 with a sugar group and at position 4 and/or position 6 with an amino acid group. The compds. are based on the structure of phosphatidylinositol hexamannosides (PIM6) of Mycobacteria and may be used as mimics of the naturally occurring PIMs in order to induce biol. responses normally attributed to the natural compd. or may be used as biol. inert carriers in order to deliver specific pharmaceutically active compds. to lipid rafts/caveolae (no data). Thus,

triethylammonium-[2-O-(.alpha.-D-mannopyranosyl)-L-myo-inosit-1-yl]-

[(2R)-2,3-bis(myristoyloxy)propyl]-phosphate was prepd. as phosphatidylinositol hexamannoside mimic and potential drug delivery agent.

IT 579493-79-1P

(prepn. of substituted inositols and their use as phosphatidylinositol hexamannoside mimics and potential drug delivery agents)

RN 579493-79-1 HCAPLUS

CN D-myo-Inositol, 3,4,5-tris-O-(phenylmethyl)-6-O-2-propenyl-2-O-

[2,3,4,6-tetrakis-O-(phenylmethyl)-.alpha.-D-mannopyranosyl]-,
(2R)-2,3-bis[(1-oxotetradecyl)oxy]propyl hydrogen phosphate (9CI) (CA
INDEX NAME)

Absolute stereochemistry.

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IPCI C07H0015-207 [ICM,7]; A61K0031-70 [ICS,7]
IPCR C07H0015-207 [I,A]
CC
     33-6 (Carbohydrates)
     Section cross-reference(s): 63
ΙT
     22006-88-8P, Liriodendritol
                                    55123-24-5P
                                                  111901-82-7P
     120202-94-0P
                    126722-28-9P
                                                   154372-19-7P
                                    131233-71-1P
     154372-20-0P
                    154459-79-7P
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     154459-83-3P
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     170900-79-5P
                    380366-30-3P
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        (prepn. of substituted inositols and their use as
        phosphatidylinositol hexamannoside mimics and potential drug
        delivery agents)
REFERENCE COUNT:
                          4
                                THERE ARE 4 CITED REFERENCES AVAILABLE FOR
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L48 ANSWER 11 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2003:456232 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 139:22337

TITLE: New phosphonate derivatives, their preparation method, and their use as modulators of the

RE FORMAT

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

activity of t.gamma.9.delta.2 lymphocytes INVENTOR(S): Montero, Jean Louis; Zgani, Ibrahim; Menut,

Chantal; Gallois, Valerie

PATENT ASSIGNEE(S): Laboratoires Mayoly Spindler, Fr.; Centre National

de la Recherche Scientifique CNRS; Universite

Montpellier II

SOURCE: Fr. Demande, 118 pp.

CODEN: FRXXBL

DOCUMENT TYPE: Patent LANGUAGE: French

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

	TENT				KIN:		DATE				ICAT					ATE
	2833				A1		2003				001-					0011211
FR	2833	266			В1		2004	1022			,					
WO	2003	0501	28		A1		2003	0619	1	WO 2	002-	FR41	90		2	0021205
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AU	2002	3003	1 1		AI		2003	J623		AU Z		3663 	1 1		2	0021205
EP	1453	840			A1		2004	908	:	EP 2		8045	96		2	0021205
	R·	ΔТ.	BE.	CH.	DE.	DK.	ES,	FR.	GB.	GR.			T.II.	NT.	SE.	MC -
				-			FI,					-		•	-	•
JP	2005			~-,			2005					5511				0021205
US	2006	0241	087		A1		2006	1026	1	US 2	005-		13		2	0051104
US	US 20100204184				A1		2010	0812						20100114		
US	US 8017596				В2		2011	0913								
	IORITY APPLN. INFO.:			.:	22		2011	0,910		FR 2		1597: 	1	-	A 2	0011211
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									1	US 2			13		A1 2	0051104

OTHER SOURCE(S): MARPAT 139:22337

ED Entered STN: 15 Jun 2003

GΙ

The invention refers to new phosphonates A-[P(:O)(O-)-O]n-BX+[I, A=C1-C50 alkylAB which may be linear, branched, cyclic, satd. or unsatd. (alkenyl or alkynyl), (un) substituted by arom. groups, functionalized by bridging ethers, carboxylic acids, esters, amides, nitriles, hydroxyls, aldehydes, ketones, halogens, amines, thiols, thioketones, episulfides, selenols, selenoketones, sulfides, sulfones, sulfoxides, or may contain one or more heterocycles; n = 1-4; X = H or a pharmaceutically acceptable cationic org. group or mineral; B = X and -P(:0) (A) O-X+ (same X and A as above)]. Compds. I, e.g., pyrophosphonate [(E)-Me2C:CHCH:CHP(O)(O-)OP(O)(O-)2](NH4+)3, and their pharmaceutical compns. modulate the proliferation of T .gamma.9.delta.2 lymphocytes, and are useful for the treatment or prevention of infectious diseases, tumors, and chronic inflammatory diseases. A claimed method for conversion of di-Et alkylphosphonates of the invention to the corresponding phosphonic diacid is by treatment with a trimethylsilyl halide to give a bis(trimethylsilyl) phosphonate which is then hydrolyzed. A method is also claimed for prepg. compds. I which has a step comprising reaction of a phosphonic diacid with di-Ph phosphate tributylammonium chloride salt to give a phosphonic anhydride which is then treated with tributylammonium orthophosphate in pyridine. Diagnostic kits contq. compds. I for modulating activity of T .gamma.9.delta.2 lymphocytes are claimed. ΙT 537696-98-3P

(prepn. of phosphonate derivs. as modulators of T .gamma.9.delta.2 lymphocyte activity and for treatment/prevention of infectious diseases, tumors, and chronic inflammatory diseases)

RN 537696-98-3 HCAPLUS

CN

Isohypophosphoric acid, [[(2-methyl-2-propenyl)oxy]methyl]-,
triammonium salt (9CI) (CA INDEX NAME)

$$\begin{array}{c} \text{CH2} \\ \text{Me-C-CH}_2 - \text{O-CH}_2 - \text{P-OPO}_3 \text{H}_2 \\ \text{OH} \end{array}$$

●3 NH3

IPCI C07F0009-40 [ICM,7]; C07F0009-38 [ICS,7]; A61K0031-662 [ICS,7];
 A61K0031-663 [ICS,7]; A61P0031-00 [ICS,7]
IPCR G01N0033-15 [I,A]; A61K0031-662 [I,A]; A61K0031-663 [I,A]; A61K0035-14
 [I,A]; A61K0035-26 [I,A]; A61P0001-00 [I,A]; A61P0001-16 [I,A];
 A61P0019-02 [I,A]; A61P0021-00 [I,A]; A61P0029-00 [I,A]; A61P0031-00
 [I,A]; A61P0031-12 [I,A]; A61P0031-18 [I,A]; A61P0031-22 [I,A];

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A61P0035-00 [I,A]; A61P0037-00 [I,A]; A61P0037-02 [I,A]; A61P0037-04
    [I,A]; A61P0043-00 [I,A]; C07F0009-38 [I,A]; C07F0009-655 [I,A];
    C12N0005-0783 [I,A]
CC
    29-7 (Organometallic and Organometalloidal Compounds)
    Section cross-reference(s): 1
ΙT
    496042-87-6P 496042-88-7P 496042-89-8P 496042-90-1P
    496042-91-2P 496042-92-3P 537696-50-7P 537696-53-0P
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    537696-81-4P 537696-82-5P 537696-83-6P 537696-84-7P
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    537696-90-5P 537696-92-7P 537696-96-1P 537696-98-3P
    537696-99-4P 537697-00-0P 537697-01-1P
        (prepn. of phosphonate derivs. as modulators of T .gamma.9.delta.2
       lymphocyte activity and for treatment/prevention of infectious
       diseases, tumors, and chronic inflammatory diseases)
OS.CITING REF COUNT: 1
                            THERE ARE 1 CAPLUS RECORDS THAT CITE THIS
                             RECORD (1 CITINGS)
REFERENCE COUNT:
                       19
                             THERE ARE 19 CITED REFERENCES AVAILABLE FOR
                             THIS RECORD. ALL CITATIONS AVAILABLE IN THE
                             RE FORMAT
L48 ANSWER 12 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN
                    2003:282375 HCAPLUS Full-text
ACCESSION NUMBER:
DOCUMENT NUMBER:
                       138:309276
TITLE:
                       Liposomes containing (ether)-lysolecithins for
                       treating Leishmanioses and other protozoan
                       diseases
INVENTOR(S):
                       Eibl, Joerg
                       Max-Planck-Gesellschaft Zur Foerderung Der
PATENT ASSIGNEE(S):
                       Wissenschaften E.V., Germany
                       PCT Int. Appl., 59 pp.
SOURCE:
                       CODEN: PIXXD2
DOCUMENT TYPE:
                       Patent
LANGUAGE:
                       German
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:
                   KIND
                                        APPLICATION NO.
    PATENT NO.
                              DATE
                                                             DATE
                                         _____
                      ____
                              _____
    WO 2003028701
                       A2
                              20030410 WO 2002-EP10650
                                                              20020927
                                                 <--
    WO 2003028701
                       А3
                              20031224
        W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH,
            CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD,
            GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ,
            LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ,
            NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ,
            TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
        RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ,
            BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK,
            EE, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, SK, TR,
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A1 20030424 DE 2001-10148066 20010928

DE 10148066

BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG

AU 2002362507 A1 20030414 AU 2002-362507 20020927 <-
PRIORITY APPLN. INFO.:

DE 2001-10148066 A 20010928 <-
WO 2002-EP10650 W 20020927 <--

OTHER SOURCE(S): MARPAT 138:309276

ED Entered STN: 11 Apr 2003

The invention relates to a novel pharmaceutical formulation contg. (ether)lysolecithin compds. in a liposomal form; further cholesterol and a neg. charged component are included in the liposomes. The pharmaceutical formulation is esp. suitable for stimulating leucopoiesis, and for treating tumor diseases and protozoan diseases, esp. Leishmanioses and amoebic diseases, acariasis and diseases caused by arthropods. Oher drugs can be added. Thus 1-O-octadecyl-2-O-methyl-glycero-3-phosphocholine (ET18OCH3) liposomes were prepd. from 47.5 .mu.M cholesterol, 7.5 .mu.M 1,2-dioleyl-sn-glycero-3-phosphoglycerin monosodium salt and 45.0 .mu.M ET18OCH3 in 200 mL chloroform by mixing, heating and evapg. the solvent. To the residue 450 g 0.25 M 1,2-propanediol were added, heated and filtered.

IT 149143-20-4

(liposomes contg. (ether)-lysolecithins for treating Leishmanioses and other protozoan diseases) $\,$

RN 149143-20-4 HCAPLUS

CN Ethanaminium, 2-[[hydroxy[3-(octadecyloxy)-2-(2-propen-1-yloxy)propoxy]phosphinyl]oxy]-N,N,N-trimethyl-, inner salt (CA INDEX NAME)

IPCI A61K0009-10 [ICM,7]; A61P0043-00 [ICS,7]

A61P0027-02 [I,A]; A61P0031-04 [I,A]; A61P0033-00 [I,A]; A61P0035-00 [I,A]; A61P0043-00 [I,A] 63-6 (Pharmaceuticals) CC Section cross-reference(s): 14 57-55-6, 1,2-Propane diol, biological studies 57-88-5, Cholesterol, ITbiological studies 64-17-5, Ethanol, biological studies 67-63-0, 2-Propanol, biological studies 78-92-2, 2-Butanol 79-57-2, Oxytetracycline 126-07-8, Griseofulvin 564-25-0, Doxycycline 1397-89-3, Amphotericin B 4358-16-1D, esters with poly glycerols 10118-90-8, Minocycline 70641-51-9 77249-78-6 77286-66-9 78858-43-2 79217-60-0, Cyclosporin 87746-72-3 91605-33-3 102340-77-2 105405-91-2 149143-20-4 508174-70-7 508174-73-0 508182-36-3 508182-37-4 508190-02-1 508190-03-2 508190-04-3 508190-05-4 508190-06-5 508190-07-6 508190-08-7508190-09-8 508190-10-1 508190-11-2 508190-12-3 508190-13-4 508190-14-5 508190-15-6 508190-16-7 508190-17-8 508190-18-9 508190-19-0 508190-20-3 508190-21-4 509076-44-2 509076-45-3

IPCR A61K0009-127 [I,A]; A61K0031-575 [I,A]; A61K0031-685 [I,A];

(liposomes contg. (ether)-lysolecithins for treating Leishmanioses and other protozoan diseases)

OS.CITING REF COUNT: 1 THERE ARE 1 CAPLUS RECORDS THAT CITE THIS

RECORD (1 CITINGS)

REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 13 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2002:794472 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 138:24910

TITLE: Syntheses and Calcium-Mobilizing Evaluations of

N1-Glycosyl-Substituted Stable Mimics of Cyclic

ADP-Ribose

AUTHOR(S): Huang, Li-Jun; Zhao, Yong-Yuan; Yuan, Lan; Min,

Ji-Mei; Zhang, Li-He

CORPORATE SOURCE: National Key Laboratory of Natural and Biomimetic

Drugs, Peking University, Beijing, 100083, Peop.

Rep. China

SOURCE: Journal of Medicinal Chemistry (2002), 45(24),

5340-5352

CODEN: JMCMAR; ISSN: 0022-2623

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal LANGUAGE: English

OTHER SOURCE(S): CASREACT 138:24910

ED Entered STN: 20 Oct 2002

AΒ Cyclic ADP-ribose (cADPR) is not only a potent endogenous calcium modulator but also a second messenger. However, studies on the mechanism of cADPR action were limited due to its instability and lack of available structural modifications in the N1-glycosyl unit of cADPR. In the present work, a series of N1-glycosyl mimics with different configurational glycosyls or an ether strand were designed and synthesized mimicking the furanose ring. SN2 substitutions were carried out between the protected inosine and glycosyl triflates to form the N1-glycosylinosine derivs., accompanied with some O6-glycosyl-substituted derivs. as side products. The intramol. cyclization followed the strategy described by Matsuda et al. It was found that the 8-unsubstituted substrate could also be used to construct the intramol. cyclic pyrophosphate. The activities of N1-glycosyl-substituted cADPR mimics were evaluated by induced Ca2+ release in rat brain microsomes and HeLa cells. It was found that the configuration of the N1-qlycosyl moiety in cADPR is not a crit. structural factor for retaining the activity of mobilizing Ca2+ release. More interestingly, the N1-acyclic analog exhibited strong activity by inducing Ca2+ release in both rat brain microsomes and HeLa cells. It constitutes a useful tool for further studies.

IT 478044-49-4P

(syntheses and calcium-mobilizing evaluations of N1-glycosyl-substituted stable mimics of cyclic ADP-ribose)

RN 478044-49-4 HCAPLUS

CN D-Arabinitol, 2,5-anhydro-4-[8-bromo-9-(2,3-di-O-acetyl-5-O-phosphono-beta.-D-ribofuranosyl)-6,9-dihydro-6-oxo-1H-purin-1-yl]-4-deoxy-3-O-2-propenyl-, 1-(S-phenyl hydrogen phosphorothioate) (9CI) (CA INDEX NAME)

Absolute stereochemistry.

33-9 (Carbohydrates) CC Section cross-reference(s): 1, 6 ΙT

444989-14-4P 444989-15-5P

444989-16-6P 444989-17-7P 444989-18-8P 478044-19-8P 478044-20-1P 478044-21-2P 478044-23-4P 478044-25-6P 478044-22-3P 478044-24-5P 478044-27-8P 478044-28-9P 478044-29-0P 478044-30-3P 478044-31-4P 478044-32-5P 478044-33-6P 478044-34-7P 478044-35-8P 478044-36-9P 478044-37-0P 478044-38-1P 478044-39-2P 478044-40-5P 478044-41-6P 478044-42-7P 478044-44-9P 478044-46-1P 478044-47-2P 478044-48-3P

478044-49-4P 478044-51-8P 478044-53-0P 478044-54-1P 478044-59-6P 478044-63-2P

(syntheses and calcium-mobilizing evaluations of

N1-glycosyl-substituted stable mimics of cyclic ADP-ribose)

OS.CITING REF COUNT: 30 THERE ARE 30 CAPLUS RECORDS THAT CITE THIS

RECORD (30 CITINGS)

REFERENCE COUNT: 34 THERE ARE 34 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 14 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2002:516677 HCAPLUS Full-text

DOCUMENT NUMBER: 137:57599

TITLE: Prevention and treatment of pulmonary bacterial

infection or symptomatic pulmonary exposure to endotoxin by inhalation of anti-endotoxin drugs

Rossignol, Daniel P.; Vermeulen, Mary W. INVENTOR(S):

PATENT ASSIGNEE(S): Eisai Co., Ltd., Japan

U.S., 37 pp., Cont.-in-part of U.S. 293,856. SOURCE:

CODEN: USXXAM

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 6417172	В1	20020709	US 1999-449601	19991123
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US 5750664	A	19980512	US 1995-461675	19950605
			<	
US 5935938	А	19990810	US 1996-658656	19960605
			<	
US 6184366	В1	20010206	US 1999-293856	19990416
55 525 500	21		00 1333 230000	

CA	2392356		A1	20010531	CA 2000-2392356	20001122
WO	20010378	43	A1	20010531		20001122
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EP	CY,	DE, DK,	ES,	FI, FR, GB, CG, CI, CM,	SL, SZ, TZ, UG, ZW, GR, IE, IT, LU, MC, GA, GN, GW, ML, MR, EP 2000-980723	NL, PT, SE, NE, SN, TD, TG
EP	1248629		В1	20050126	<	
					GB, GR, IT, LI, LU, MK, CY, AL, TR	NL, SE, MC,
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AT	287719		Т	20050215	AT 2000-980723	20001122
ES	2237475		Т3	20050801	ES 2000-980723	20001122
US	20030134	805	A1	20030717	US 2002-167222	20020611
	6683063 1051490		B2 A1	20040127 20050422	нк 2003-102773 <	20030416
JP	20072698	12	A	20071018	JP 2007-154127	20070611
	4712001 20111219	70	B2 A	20110629 20110623	JP 2011-18401 <	20110131
PRIORITY	Y APPLN.	INFO.:				A2 19950605
						A1 19960605
					US 1999-293856	A2 19990416
					< JP 1997-501868	A3 19960605
					< US 1999-449601	A 19991123
					< WO 2000-US32177	W 20001122
					< JP 2007-154127	A3 20070611

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 137:57599

ED Entered STN: 11 Jul 2002

GΙ

Disaccharide compds. I, wherein R is H, CH2OH, alkoxide; R1 is acyl; R2 is C5-C15 alkyl R3 is C5-C18 alkyl, acyl, R4 is C4-C20 alkyl, oxyalkyl; A1 and A2 are independently OH, phosphate, phosphonate, ester; were prepd. for and treatment of pulmonary bacterial infection or symptomatic pulmonary exposure to endotoxin. The invention provides methods of preventing and treating pulmonary bacterial infection or symptomatic pulmonary exposure to endotoxin and related conditions in a patient by administering to the patient anti-endotoxin compds. by inhalation. The invention provides methods of preventing and treating pulmonary bacterial infection or symptomatic pulmonary exposure to endotoxin and related conditions in a patient by administering to the patient anti-endotoxin compds. by inhalation. Thus, disaccharide lipid II was prepd. and tested in mice and suppressed the prodn. of TNF following administration of LPS.

IT 234088-16-5P

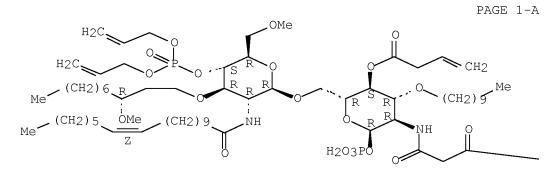
(prevention and treatment of pulmonary bacterial infection or symptomatic pulmonary exposure to endotoxin by inhalation of anti-endotoxin drugs such as disaccharide lipid A analogs in relation to inhibition of cytokine prodn.)

RN 234088-16-5 HCAPLUS

CN .alpha.-D-Glucopyranose, 6-O-[4-O-[bis(2-propenyloxy)phosphinyl]-2-deoxy-3-O-[(3R)-3-methoxydecyl]-6-O-methyl-2-[[(11Z)-1-oxo-11-octadecenyl]amino]-.beta.-D-glucopyranosyl]-3-O-decyl-2-deoxy-2-[(1,3-dioxotetradecyl)amino]-, 4-(3-butenoate) 1-(dihydrogen phosphate), disodium salt (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Double bond geometry as shown.



●2 Na

PAGE 1-B

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INCL 514053000
IPCI A61K0031-70 [ICM, 7]
IPCR C07H0003-04 [I,A]; C07H0011-00 [I,A]; C07H0015-04 [I,A]; C07H0015-12
     [I,A]
NCL
    514/053.000
     1-12 (Pharmacology)
CC
     Section cross-reference(s): 26, 33, 63
                               36875-26-0P
                                             128313-03-1P 128387-27-9P
ΙT
     5509-08-0P
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     137766-83-7P
                    139629-59-7P
                                   139686-99-0P
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     185954-75-0P
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                    185955-13-9P
                                   185955-14-0P
                                                   185955-15-1P
     185955-16-2P
                    185955-17-3P
                                   185955-18-4P
                                                   185955-19-5P
                    185955-21-9P
     185955-20-8P
                                   185955-22-0P
                                                   185955-23-1P
     185955-24-2P
                    185955-25-3P
                                   185955-26-4P
                                                   185955-28-6P
     185955-29-7P
                    234088-12-1P
                                   234088-13-2P
                                                   234088-14-3P
     234088-15-4P
                    234088-16-5P
                                   234088-19-8P
                                                   234088-20-1P
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10/596.747

234088-21-2P 234088-22-3P 234088-23-4P 234088-24-5P (prevention and treatment of pulmonary bacterial infection or symptomatic pulmonary exposure to endotoxin by inhalation of

anti-endotoxin drugs such as disaccharide lipid A analogs in relation to inhibition of cytokine prodn.)

OS.CITING REF COUNT: 3 THERE ARE 3 CAPLUS RECORDS THAT CITE THIS

RECORD (3 CITINGS)

REFERENCE COUNT: 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 15 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2002:107823 HCAPLUS Full-text

DOCUMENT NUMBER: 136:156515

TITLE: Adhesive compositions containing monophosphates

for bonding to hard tissues

INVENTOR(S): Stangel, Ivan; Xu, Jingwei; Ellis, Thomas; Sacher,

Edward

PATENT ASSIGNEE(S): Biomat Sciences, Inc., USA

SOURCE: U.S. Pat. Appl. Publ., 5 pp., Cont.-in-part of

Appl. No. PCT/US99/18582.

CODEN: USXXCO

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 20020015682	A1	20020207	US 2001-785555 <	20010220
US 6645952	В2	20031111		
WO 2000010478	A1	20000302	WO 1999-US18582	19990817
			<	
W: AU, CA, C	N, JP, US	3		
RW: AT, BE, C	H, CY, DE	E, DK, ES,	FI, FR, GB, GR, IE,	IT, LU, MC,
NL, PT, S	E			
US 20030118523	A1	20030626	US 2003-339621	20030110
			<	
US 6664245	В2	20031216		
PRIORITY APPLN. INFO.:			US 1998-96838P	P 19980818
			<	
			WO 1999-US18582	A2 19990817
			<	

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

ED Entered STN: 10 Feb 2002

AB An adhesive compn. is provided which is suitable for the bonding of polymeric materials, in whole or in part, such as composite resins, or methacrylate-contg. glass-ionomer filling materials, to tooth enamel, or tooth dentin, or to other hand tissues of the human body, such as bone. The compn. comprises an unsatd. carboxylic acid ester, an unsatd. phosphate ester and other crosslinking agents. The esp. preferred ethylenically unsatd. monophosphates, (CH2::C(R)CH2OCH2)nR1OP(O)(OH)2 (R = H, C1-4 alkyl, CN; R1 = aliph., cycloaliph., aryl), are provided as new compds. For example, a mixt. of 2-hydroxyethyl methacrylate (HEMA) (35% by wt.) and

US 2001-785555

<--

A3 20010220

pentaerythritol triallyl ether monophosphate acid ester (PTEPAE) (10% by wt.), the balance being acetone, was prepd. and the bond strength of the soln. to dentin was tested. The mean peak stress at failure for seven dentin samples was 20.5 MPa. However, the mean peak stress at failure for six dentin samples etched by 35% phosphoric acid was 25.5 MPa.

IT 259250-33-4

(adhesive compns. for bonding of polymeric materials to hard tissues)

RN 259250-33-4 HCAPLUS

CN 1-Propanol, 3-(2-propen-1-yloxy)-2,2-bis[(2-propen-1-yloxy)methyl]-, 1-(dihydrogen phosphate) (CA INDEX NAME)

INCL 424049000

IPCI A61K0007-16 [ICM]; C07F0009-113 [ICS]
IPCR A61K0006-00 [I,A]; C07F0009-09 [I,A]

NCL 424/049.000; 558/208.000; 514/112.000; 514/129.000; 558/183.000

CC 63-7 (Pharmaceuticals)

Section cross-reference(s): 23

IT 259250-33-4 395068-31-2

(adhesive compns. for bonding of polymeric materials to hard tissues)

L48 ANSWER 16 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2001:693328 HCAPLUS Full-text

DOCUMENT NUMBER: 135:257469

DOCUMENT NUMBER: 133:23/409

TITLE: Preparation of caprolactam derivatives with Src-SH2 domain inhibitor activity and their

intermediates, and their application as bone

resorption inhibitors

INVENTOR(S): Deprez, Pierre; Lesuisse, Dominique; Benard,

Didier

PATENT ASSIGNEE(S): Ariad Pharmaceuticals, Inc., USA

SOURCE: PCT Int. Appl., 43 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2001068655	A2	20010920	WO 2001-US7935 <	20010312

WO 2001068655 A3 20020801

W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, HU, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW,

MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZA RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG AU 2001043601 Α 20010924 AU 2001-43601 20010312 <--PRIORITY APPLN. INFO.: US 2000-523243 A 20000310 <--WO 2001-US7935 W 20010312 <--

OTHER SOURCE(S): MARPAT 135:257469

ED Entered STN: 21 Sep 2001

GΙ

Caprolactam derivs. [I; X = XH2, N; R1 = alkyl, alkenyl, alkynyl, arylalkyl, aryl-alkenyl, aryl-alkynyl, (un)substituted cycloalkyl etc.; A1 = CH(Z)-alkylaryl, CH(Z):CH-alkylaryl, CH(Z)-aryl, alkylaryl, aryl; Z = H, tetrazole, (un)substituted NH2 and CONH2; A2 = P(O)(OH)2 or esters, B(OH)2 or esters, various carboxylic or sulfonic acids or their derivs.], as well as isomers, physiol. acceptable salts, and/or prodrugs, were prepd. as inhibitors of the Src-SH2 receptor. Thus, caprolactam deriv. (II) was prepd. via a multistep synthetic sequence starting from (2S)-N-Boc-Tyr-O-(PO3Bn2)and (S)-3-amino-hexahydro-2H-azepin-2-one. In a scintillation proximity assay for inhibition of the binding of the ligand [125I]-EPQpYEEIPIYL to biotinylated SH2 protein, II had an IC50 of 0.009 .mu.M, vs. 0.2-0.4 .mu.M for the ref. peptide PYEEI. IT 361385-61-7DP, Wang resin-bound

(prepn. of caprolactam derivs. as Src-SH2 domain antagonists)

RN 361385-61-7 HCAPLUS

CN D-Tyrosine, N-acetyl-, 2-propenyl ester, 1,1-dimethylethyl hydrogen phosphate (ester) (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IPCI C07F0009-02 [ICM, 7]

IPCR A61K0038-00 [N,A]; C07F0009-12 [I,A]; C07F0009-40 [I,A]; C07F0009-553

[I,A]; C07K0005-06 [I,A]; C07K0005-065 [I,A]
34-2 (Amino Acids, Peptides, and Proteins)

CC 34-2 (Amino Acids, Peptides, and Pro Section cross-reference(s): 1, 27

IT 76944-95-1P 361385-59-3P 361385-60-6P 361385-61-7DP,

Wang resin-bound 361385-62-8DP, Wang resin-bound 361385-63-9P

361385-64-0P 361385-65-1P

(prepn. of caprolactam derivs. as Src-SH2 domain antagonists)

OS.CITING REF COUNT: 6 THERE ARE 6 CAPLUS RECORDS THAT CITE THIS

RECORD (6 CITINGS)

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 17 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2001:396678 HCAPLUS Full-text

DOCUMENT NUMBER: 135:528

TITLE: Prevention and treatment of pulmonary bacterial

infection or symptomatic pulmonary exposure to endotoxin by inhalation of antiendotoxin drugs

INVENTOR(S): Rossignol, Daniel P.; Vermeulen, Mary W.

PATENT ASSIGNEE(S): Eisai Co., Ltd., Japan SOURCE: PCT Int. Appl., 87 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 4

PATENT INFORMATION:

PAT	CENT 1	NO.			KIN	D	DATE			APPL	ICAT	ION :	NO.		D	ATE	
WO	2001	0378	43		A1	_	2001		1	WO 2		 US32 	177		2	0001	
	W: RW:	CN, GM, LR, PL, UA, GH,	CR, HR, LS, PT, UG, GM,	CU, HU, LT, RO, US, KE,	CZ, ID, LU, RU, UZ, LS,	DE, IL, LV, SD, VN, MW,	AU, DK, IN, MA, SE, YU, MZ, FR,	DM, IS, MD, SG, ZA, SD,	DZ, JP, MG, SI, ZW SL,	EE, KE, MK, SK,	ES, KG, MN, SL,	FI, KP, MW, TJ,	GB, KR, MX, TM,	GD, KZ, MZ, TR,	GE, LC, NO, TT,	GH, LK, NZ, TZ,	
	6417 2392	TR,	BF,	BJ,	CF,	CG,	CI, 2002	CM, 0709	GA,	GN, US 1	GW, 999- <	ML, 4496 	MR, 01	NE,	SN,	•	123
											<						

EP	1248629			A1	20021	016	EP	2000-9			20001122
EP	1248629			B1			an ar	·			
					DK, ES, I	•	· ·	-		NL, SI	±, MC,
JP	20035148		•	•		•	•	2001-5	39457		20001122
AT	287719			Т	20050	215	АТ	2000-9	80723		20001122
HK	1051490			A1	20050	422	HK	2003-1	02773		20030416
PRIORITY	APPLN.	INFO	.:				US		49601	A2	19991123
							US	1995-4	61675 -	A2	19950605
							US		58656	A1	19960605
							US	,	93856	A2	19990416
							WO	,	s32177	M	20001122

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT OTHER SOURCE(S): MARPAT 135:528

ED Entered STN: 01 Jun 2001

AB The invention provides methods of preventing and treating pulmonary bacterial infection or symptomatic pulmonary exposure to endotoxin and related conditions in a patient by administering to the patient antiendotoxin compds. by inhalation.

IT 234088~16~5P

(prevention and treatment of pulmonary bacterial infection or symptomatic pulmonary exposure to endotoxin by inhalation of antiendotoxin drugs such as lipid A analogs in relation to inhibition of cytokine prodn.)

RN 234088-16-5 HCAPLUS

CN .alpha.-D-Glucopyranose, 6-O-[4-O-[bis(2-propenyloxy)phosphinyl]-2-deoxy-3-O-[(3R)-3-methoxydecyl]-6-O-methyl-2-[[(11Z)-1-oxo-11-octadecenyl]amino]-.beta.-D-glucopyranosyl]-3-O-decyl-2-deoxy-2-[(1,3-dioxotetradecyl)amino]-, 4-(3-butenoate) 1-(dihydrogen phosphate), disodium salt (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry as shown.

●2 Na

PAGE 1-B

IPCI A61K0031-70 [ICM, 7] IPCR A61K0009-12 [I,A]; A61K0009-00 [I,A]; A61K0031-6615 [I,A]; A61K0031-70 [I,A]; A61K0045-00 [I,A]; A61P0011-00 [I,A]; A61P0031-00 [I,A] CC1-12 (Pharmacology) Section cross-reference(s): 26, 33, 63 36875-26-0P ΙT 5509-08-0P 19525-80-5P 128313-03-1P 128387-27-9P 137766-83-7P 139629-59-7P 139686-99-0P 185954-74-9P 185954-75-0P 185954-76-1P 185954-77-2P 185954-78-3P 185954-79-4P 185954-80-7P 185954-81-8P 185954-82-9P 185954-83-0P 185954-84-1P 185954-85-2P 185954-86-3P 185954-87-4P 185954-88-5P 185954-89-6P 185954-90-9P 185955-12-8P 185955-13-9P 185955-14-0P 185955-15-1P 185955-16-2P 185955-17-3P 185955-18-4P 185955-19-5P 185955-20-8P 185955-21-9P 185955-22-0P 185955-23-1P 185955-24-2P 185955-25-3P 185955-26-4P 185955-28-6P 185955-29-7P 234088-12-1P 234088-13-2P 234088-14-3P 234088-15-4P 234088-16-5P 234088-19-8P 234088-20-1P 234088-21-2P 234088-22-3P 234088-23-4P 234088-24-5P (prevention and treatment of pulmonary bacterial infection or

symptomatic pulmonary exposure to endotoxin by inhalation of antiendotoxin drugs such as lipid A analogs in relation to inhibition of cytokine prodn.)

REFERENCE COUNT: 1 THERE ARE 1 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 18 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN 2000:144683 HCAPLUS Full-text ACCESSION NUMBER:

DOCUMENT NUMBER: 132:185486

TITLE: Adhesive compositions containing ethylenically

unsaturated monophosphates for the hard tissues of the human body such as tooth enamel and tooth

dentin

Xu, Jingwei; Stangel, Ivan; Ellis, Thomas; Sacher, INVENTOR(S):

Edward

PATENT ASSIGNEE(S): Biomat Services, Inc., USA

SOURCE: PCT Int. Appl., 14 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

	PATENT NO.							APPLICATION NO.								
										WO		US185				19990817
			AT,		CH,	JP, CY,		, DK,	ES,	FI, F	R, GB,	GR,	IE,	IT,	LU	J, MC,
	CA	2340				A1		2000	0302		1999- <		900			19990817
	AU	9954	876			А		2000	0314	AU		54876	6			19990817
	ΕP	1105	063			A1		2001	0613			94116	68			19990817
		R:	-	BE, IE,		DE,	DK,	, ES,	FR,	GB, G			LU,	NL,	SE	, MC,
	JP	2002	•			Т		2002	0730		2000-		04			19990817
	CN	1003	8973	1		С		2008	0528	CN	1999-		65			19990817
	US	2002	0015	682		A1		2002	0207	US	2001-		55			20010220
	US	6645	952			В2		2003	1111							
	US	2003	0118	523		A1		2003	0626	US	2003-		21			20030110
		6664				В2		2003	1216							
PRIO:	RIT	Y APP	LN.	INFO	.:					US	1998-				Р	19980818
											1999-			1	W	19990817
										US	2001-	7855	55		АЗ	20010220

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

10/596.747

OTHER SOURCE(S): MARPAT 132:185486

ED Entered STN: 03 Mar 2000

AB Title compns. are suitable for the bonding of polymeric materials, in whole or in part, such as composite resins, or methacrylate-contg. glass-ionomer filling materials, to tooth enamel, or tooth dentin, or to other hard tissues of the human body, such as bone. The compns. comprise unsatd. carboxylic acid esters, unsatd. phosphate esters represented by the formula [CH2:C(R1)CH2OCH2]nROPO3H2, and other crosslinking agents, where R1 is a hydrogen atom, C1-C4 alkyl, or CN; R is an aliph., cycloaliph., or aryl radical contg. from 1 to 10 carbon atoms and having a valence of n + 1; and n is an integer from 1 to 5.

IT 259250-34-5P

(prepn. of dental adhesive compns. contg. ethylenically unsatd. monophosphates)

RN 259250-34-5 HCAPLUS

CN 2-Propenoic acid, 2-methyl-, 2-hydroxyethyl ester, polymer with 3-(2-propenyloxy)-2,2-bis[(2-propenyloxy)methyl]propyl dihydrogen phosphate (9CI) (CA INDEX NAME)

CM 1

CRN 259250-33-4 CMF C14 H25 O7 P

$$CH_2 - OPO_3H_2$$
 $H_2C = CH - CH_2 - O - CH_2 - CH_2 - CH_2 - CH_2 - CH_2 - CH_2 - CH_2$
 $CH_2 - O - CH_2 - CH_2 - CH_2$
 $CH_2 - O - CH_2 - CH_2 - CH_2$

CM 2

CRN 868-77-9 CMF C6 H10 O3

$$\begin{array}{ccc} ^{\rm H2C} & {\rm O} \\ \parallel & \parallel \\ {\rm Me-C-C-O-CH_2-CH_2-OH} \end{array}$$

CC 63-7 (Pharmaceuticals)

IT 259250-34-5P 259250-35-6P

(prepn. of dental adhesive compns. contg. ethylenically unsatd. monophosphates)

IT 259250-33-4P

(prepn. of ethylenically satd. monophosphates for dental adhesives)
OS.CITING REF COUNT: 7 THERE ARE 7 CAPLUS RECORDS THAT CITE THIS
RECORD (7 CITINGS)

REFERENCE COUNT: 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 19 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 2000:117065 HCAPLUS Full-text

DOCUMENT NUMBER: 132:161274

TITLE: CMP derivative sialyltransferase inhibitors,

preparation, pharmaceutical compositions, and

WO 1999-EP5697

<--

W 19990806

therapeutic use

INVENTOR(S): Schmidt, Richard R.; Schaub, Christoph; Muller,

Bernd; Amann, Franz

PATENT ASSIGNEE(S): Germany

SOURCE: PCT Int. Appl., 63 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.					KIN	D	DATE			APPL	ICAT	ION 1	NO.		Di	ATE
WO 2	WO 2000008040				A1 20000217			WO 1999-EP5697 <						1	9990806	
	W:	CZ,	DE,	DK,	EE,	ES,	AZ, FI,	GB,	GD,	GE,	BR, GH,	BY, GM,	HR,	HU,	ID,	IL,
		MD,	MG,	MK,	MN,	MW,	KP, MX, TR,	NO,	NZ,	PL,	PT,	RO,	RU,	SD,	SE,	SG,
	RW:	GH,	GM,	KE,	LS,	MW,	SD, GR,	SL,	SZ,	UG,	ZW,	AT,	BE,	CH,	CY,	DE,
AU S	9960 ⁻	•	•	•	CM, A	•	GN, 2000	•	•	•	•	•	•	TG	1	9990806
PRIORITY APPLN. INFO.:											,		OP		P 1	9980807

OTHER SOURCE(S): MARPAT 132:161274

ED Entered STN: 18 Feb 2000

GΙ

AB The invention provides potent inhibitors of sialyltransferases. The sialyltransferase inhibitors are useful for inhibiting the synthesis of sialylated

glycosides. The sialyltransferase inhibitors find use in the modulation of biol. processes that involve sialyl glycoside-mediated cell adhesion. The compds. of the invention may be used e.g. to treat inflammation.

IT 218939-23-2

(reaction; CMP deriv. sialyltransferase inhibitors, prepn., pharmaceutical compns., and therapeutic use)

RN 218939-23-2 HCAPLUS

CN 5'-Cytidylic acid, N-acetyl-, 2',3'-diacetate, monoester with 5-(acetylamino)-2,6-anhydro-1-C-[bis(2-propenyloxy)phosphinyl]-3,5-dideoxy-D-glycero-D-galacto-non-2-enitol 4,7,8,9-tetraacetate, compd. with N,N-diethylethanamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 218939-22-1 CMF C40 H54 N4 O24 P2

Absolute stereochemistry.

CM 2

CRN 121-44-8 CMF C6 H15 N

Et | Et-N-Et

REFERENCE COUNT: 8 THERE ARE 8 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 20 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1999:505657 HCAPLUS $\underline{\text{Full-text}}$

DOCUMENT NUMBER: 131:130224

TITLE: Substituted liposaccharides useful in the

treatment and prevention of endotoxemia

INVENTOR(S): Christ, William J.; Rossignol, Daniel P.;

Kobayashi, Seiichi; Kawata, Tsutomu

PATENT ASSIGNEE(S): Eisai Co., Ltd., Japan

SOURCE: U.S., 40 pp.

CODEN: USXXAM

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 4

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 5935938	 A	19990810	US 1996-658656 <	19960605
US 5681824	А	19971028	US 1995-461677	19950605
US 5750664	А	19980512	US 1995-461675	19950605
ZA 9604666	А	19970311	ZA 1996-4666 <	19960605
CN 1192216	А	19980902	CN 1996-195890	19960605
CN 1067082	С	20010613		
PT 853627	E	20040531	PT 1996-923234 <	19960605
ES 2214543	Т3	20040916	ES 1996-923234	19960605
US 6184366	В1	20010206	US 1999-293856 <	19990416
US 6417172	B1	20020709	US 1999-449601	19991123
US 20020028927	A1	20020307	US 2001-774541	20010130
US 20030144503	A1	20030731	US 2002-144670	20020513
US 20030134805	A1	20030717	US 2002-167222	20020611
US 6683063	В2	20040127		
JP 2007269812	А	20071018	JP 2007-154127 <	20070611
JP 4712001	В2	20110629		
US 20080214802	A1	20080904	US 2007-830412 <	20070730
US 7737129	В2	20100615		
US 20100227835	A1	20100909	US 2010-781166 <	20100517
US 7994154	B2	20110809		

JP 2011121970	A	20110623	JP	2011-18401		20110131
PRIORITY APPLN. INFO.:			US	1995-461675	A2	19950605
			JP	1997-501868	A3	19960605
			US	1996-658656	A1	19960605
			US	1999-293856	Α2	19990416
			US	1999-449601	A1	19991123
			US	2001-774541	В1	20010130
			US	2002-144670	В1	20020513
			JP	2007-154127	A3	20070611
			US	2007-830412	A1	20070730

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 131:130224

ED Entered STN: 16 Aug 1999

GΙ

- Novel substituted liposaccharides I (Rl = acyl; R2 = C5 to C15 alkyl; R3 = C5 to C18 acyl-alkenyl or acyl-alkynyl; R4 = C4 to C20 alkoxy-substituted alkyl; RA = CH2O-X where X is H or alkyl group; A1,A2 = OH, PO4H2, O-alkyl-OPO3H2, etc.) useful as in the prophylactic and affirmative treatment of endotoxemia including sepsis, septicemia and various forms of septic shock are prepd. Also provided are processes for prepg. the compds., e.g. II, and intermediates useful therein. The aminodeoxy disaccharide analogs inhibit tumor necrosis factor prodn. in vivo, exhibiting IC5Os between 1.5 nM and 159 nM.
- IT 234088-16-5P

(prepn. of substituted lipodisaccharides useful in the treatment and prevention of endotoxemia) $\,$

RN 234088-16-5 HCAPLUS

CN .alpha.-D-Glucopyranose, 6-O-[4-O-[bis(2-propenyloxy)phosphinyl]-2-deoxy-3-O-[(3R)-3-methoxydecyl]-6-O-methyl-2-[(11Z)-1-oxo-11-octadecenyl]amino]-.beta.-D-glucopyranosyl]-3-O-decyl-2-deoxy-2-[(1,3-dioxotetradecyl)amino]-, 4-(3-butenoate) 1-(dihydrogen phosphate), disodium salt (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Double bond geometry as shown.

^{*} STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

●2 Na

PAGE 1-B

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INCL 514053000
IPCI A61K0031-70 [ICM, 6]; C13K0005-00 [ICS, 6]
IPCR C07H0015-12 [I,A]; A61K0031-70 [I,A]; A61K0031-7028 [I,A]; A61P0031-00
     [I,A]; A61P0031-04 [I,A]; C07H0003-04 [I,A]; C07H0011-00 [I,A];
     C07H0011-04 [I,A]; C07H0013-06 [I,A]; C07H0015-04 [I,A]
NCL 514/053.000; 536/017.200; 536/017.900; 536/123.130
CC
     33-7 (Carbohydrates)
     Section cross-reference(s): 1
ΙT
     19525-80-5P
                   36875-26-0P
                                 41233-29-8P
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     234088-12-1P
                    234088-13-2P
                                   234088-14-3P
                                                   234088-15-4P
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234088-16-5P 234088-17-6P 234088-18-7P 234088-19-8P 234088-20-1P 234088-21-2P 234088-22-3P 234088-23-4P 234088-24-5P

(prepn. of substituted lipodisaccharides useful in the treatment and prevention of endotoxemia)

OS.CITING REF COUNT: 28 THERE ARE 28 CAPLUS RECORDS THAT CITE THIS

RECORD (28 CITINGS)

REFERENCE COUNT: 17 THERE ARE 17 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 21 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1999:499942 HCAPLUS Full-text

DOCUMENT NUMBER: 131:257780

TITLE: 5'-Phosphoramidates and 5'-Diphosphates of 2'-O-Allyl-.beta.-D-arabinofuranosyl-uracil,

-cytosine, and -adenine: Inhibition of

Ribonucleotide Reductase

AUTHOR(S): Manfredini, Stefano; Baraldi, Pier Giovanni;

Durini, Elisa; Vertuani, Silvia; Balzarini, Jan;

De Clercq, Erik; Karlsson, Anna; Buzzoni,

Valentina; Thelander, Lars

CORPORATE SOURCE: Department of Pharmaceutical Sciences, Ferrara

University, Italy

SOURCE: Journal of Medicinal Chemistry (1999), 42(17),

3243-3250

CODEN: JMCMAR; ISSN: 0022-2623

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal LANGUAGE: English ED Entered STN: 12 Aug 1999

AB Continuing our studies on ribonucleotide reductase (RNR) mechanism-based inhibitors, we have now prepd. the diphosphates (DP) of

2'-O-allyl-1-.beta.-D-arabinofuranosyl-uracil and -cytosine and

2'-O-allyl-9-.beta.-D-arabinofuranosyl-adenine and evaluated their inhibitory activity against recombinant murine RNR. 2'-O-Allyl-araUDP proved to be inhibitory to RNR at an IC50 of 100 .mu.M, whereas 2'-O-allyl-araCDP was only marginally active (IC50 1 mM) and 2'-O-allyl-araADP was completely inactive. The susceptibility of the parent nucleosides to phosphorylation by thymidine kinase and 2'-deoxycytidine kinase was also investigated, and all nucleosides proved to be poor substrates for the above-cited kinases. Moreover, prodrugs of

2'-O-allyl-araU and -araC monophosphates, namely

2'-O-allyl-5'-(phenylethoxy-L-alanyl phosphate)-araU and -araC, were prepd. and tested against tumor cell proliferation but proved to be inactive. A mol. modeling study has been conducted in order to explain our results. The data confirm that for both the natural and analog nucleoside diphosphates, the principal determinant interaction with the active site of RNR is with the diphosphate group, which forms strong hydrogen bonds with Glu623, Thr624, Ser625, and Thr209. Our findings indicate that the poor phosphorylation may represent an explanation for the lack of marked in vitro cytostatic activity of the test compds.

IT 245078-03-9P

(prepn. and ribonucleotide reductase inhibition of phosphoramidates and diphosphates of allyl-D-arabinofuranosyl-uracil, -cytosine and -adenine)

RN 245078-03-9 HCAPLUS

CN 2(1H)-Pyrimidinone, 4-amino-1-[5-0-[hydroxy(phosphonooxy)phosphinyl]-2-

O-2-propenyl-.beta.-D-arabinofuranosyl]-, trisodium salt (9CI) (CA INDEX NAME)

Absolute stereochemistry.

●3 Na

CC 33-9 (Carbohydrates)

Section cross-reference(s): 1

IT 245078-03-9P 245078-08-4P 245078-13-1P

(prepn. and ribonucleotide reductase inhibition of phosphoramidates and diphosphates of allyl-D-arabinofuranosyl-uracil, -cytosine and -adenine)

OS.CITING REF COUNT: 18 THERE ARE 18 CAPLUS RECORDS THAT CITE THIS

RECORD (18 CITINGS)

REFERENCE COUNT: 22 THERE ARE 22 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 22 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1999:448710 HCAPLUS Full-text

DOCUMENT NUMBER: 131:170568

TITLE: Synthesis, cytostatic activity and inhibition of

ribonucleotide reductase by 5'-phosphoramidates

and 5'-diphosphates, of

2'-O-allyl-arabinofuranosyl nucleosides
AUTHOR(S): Manfredini, S.; Baraldi, P. G.; Durini, E.;

Balzarini, J.; De Clercq, E.; Karlsson, A.;

Buzzoni, V.; Thelander, L.

CORPORATE SOURCE: Department of Pharmaceutical Sciences, Ferrara

University, Italy

SOURCE: Nucleosides & Nucleotides (1999), 18(4 & 5),

1007-1008

CODEN: NUNUD5; ISSN: 0732-8311

PUBLISHER: Marcel Dekker, Inc.

DOCUMENT TYPE: Journal LANGUAGE: English ED Entered STN: 22 Jul 1999

AB A symposium reporting that the diphosphates of a series of 2'-O-allyl-1-.beta.-D-arabinofuranosyl derivs., previously obtained by the authors, have been prepd. and tested for their inhibitory activity in an in vitro

assay using R1 and R2 subunits of the purified recombinant mouse ribonucleotide reductase (RNR). 2'-O-Allyl-araU diphosphate proved to be inhibitory, with an

IC50 of 100 .mu.M. The 5'-phosphoramidate pronucleotide of 2'-O-allyl-araU was also prepd. and tested for inhibition of tumor cell proliferation.

IT 239117-77-2P

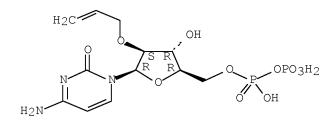
(prepn., cytostatic activity and inhibition of ribonucleotide reductase by phosphoramidates and diphosphates of

allylarabinofuranosyl nucleosides)

RN 239117-77-2 HCAPLUS

CN 2(1H)-Pyrimidinone, 4-amino-1-[5-O-[hydroxy(phosphonooxy)phosphinyl]-2-O-2-propen-1-yl-.beta.-D-arabinofuranosyl]- (CA INDEX NAME)

Absolute stereochemistry.



CC 33-9 (Carbohydrates)

Section cross-reference(s): 1

IT 239117-77-2P 239117-78-3P 239117-79-4P

239117-80-7P

(prepn., cytostatic activity and inhibition of ribonucleotide reductase by phosphoramidates and diphosphates of

allylarabinofuranosyl nucleosides)

REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 23 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1999:166753 HCAPLUS Full-text

DOCUMENT NUMBER: 130:232458

TITLE: Drug discovery using multiple membrane mimetic

affinities

INVENTOR(S): Pidgeon, Charles; Liu, Hanlan; Hauer, Kimberly;

Yin, Jiaming; Cai, Song J.

PATENT ASSIGNEE(S): Purdue Research Foundation, USA

SOURCE: PCT Int. Appl., 65 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, GM, HR, HU, ID, IL, IS, JP,

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             MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL,
             TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW
         RW: GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, CY, DE, DK,
             ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF,
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PRIORITY APPLN. INFO.:
                                             US 1997-56833P
                                                                 P 19970822
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                                             WO 1998-US17398
                                                                 W 19980821
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ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

ED Entered STN: 15 Mar 1999

AΒ The measurement of multiple membrane affinities of test compds., methods and compns. useful for acquiring data characteristic of such affinities, and a method and system for using such data alone or in combination with other mol. descriptors for the prediction of biol. activity are described. The numerical values characteristic of biol. relevant interaction of test compds. with membrane mimetic surfaces are compared with corresponding values of one or more control compds. having a known biol. activity. Probable biol. activity of a test compd. is identified with those control compds. whose multiple membrane interaction values binding data are obtained for test compds. and control compds. for use in accordance with this invention using immobilized artificial membrane chromatog. substrates in high pressure liq. chromatog. systems using aq. mobile phases. Data relevant to the thermodn. and kinetics of compd./membrane interaction is reflected in retention time and peak width, resp. All data are preferably normalized relative to a std. compd. or a set of compds., for example, a set of compds. having a common biol. activity or function. This invention also provides novel carboxylfunctional, head group-protected phospholipids useful for prepg. immobilized artificial membrane structures useful for acquiring membrane interaction data. They are prepd. by novel high yielding transphosphatidylation of phosphatidylcholine derivs. using phospholipase D in the presence of protected alcs.

IT 221108-36-7P

(intermediate; drug discovery using multiple membrane mimetic affinities and use of HPLC and prepn. of protected phospholipids useful for prepn. of immobilized artificial membranes)

RN 221108-36-7 HCAPLUS

CN L-Serine, N-[(1,1-dimethylethoxy)carbonyl]-, 2-propenyl ester,

(2R)-2-[(13-carboxy-1-oxotridecyl)oxy]-3-[(1-oxotetradecyl)oxy]propyl hydrogen phosphate (ester) (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IPCI C12Q0001-00 [ICM,6]; G01N0033-53 [ICS,6]; G01N0033-566 [ICS,6]

IPCR G01N0033-50 [I,A]; B01J0020-281 [I,A]; C12Q0001-00 [I,A]; G01N0030-02
 [N,A]; G01N0030-88 [I,A]; G01N0030-89 [N,A]; G01N0033-15 [I,A];
 G01N0033-543 [I,A]

CC 1-1 (Pharmacology)

Section cross-reference(s): 26

 $\begin{tabular}{llll} $123-78-4P$ & $20559-16-4P$ & $85483-04-1P$, Oxacyclotridecane-2,13-dione \\ \end{tabular}$

88224-08-2P 102308-32-7P 108149-60-6P 115464-01-2P 116467-63-1P 117487-53-3P 119766-79-9P 143966-57-8P 203439-80-9P 221108-33-4P 221108-34-5P 221108-35-6P 221108-36-7P

(intermediate; drug discovery using multiple membrane mimetic affinities and use of HPLC and prepn. of protected phospholipids useful for prepn. of immobilized artificial membranes)

IT 221108-33-4DP, silica propylamine-immobilized 221108-34-5DP, silica propylamine-immobilized 221108-35-6DP, silica propylamine-immobilized 221108-36-7DP, silica propylamine-immobilized

(on artificial membranes; drug discovery using multiple membrane mimetic affinities and use of HPLC and prepn. of protected phospholipids useful for prepn. of immobilized artificial membranes)

OS.CITING REF COUNT: 3 THERE ARE 3 CAPLUS RECORDS THAT CITE THIS

RECORD (3 CITINGS)

REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 24 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1998:453103 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 129:216434

ORIGINAL REFERENCE NO.: 129:43999a,44002a

TITLE: Synthesis and pharmacological evaluation of a new

class of bicyclic phospholipids, designed as

platelet activating factor antagonists

AUTHOR(S): Pecanha, Emerson Poley; Fraga, Carlos Alberto

Manssour; Mauricio, Carlos; De Sant' Anna,

Rabello; De Miranda, Ana Luisa Palhares; Barreiro,

Eliezer Jesus

CORPORATE SOURCE: Laboratorio de Avaliacao e Sintese de Substancias

Bioativas (LASSBio), Faculdade de Farmacia,

Universidade Federal do Rio de Janeiro, Rio de

Janeiro, 21944-970, Brazil Farmaco (1998), 53(5), 327-336

SOURCE: Farmaco (1998), 53(5), 327-336 CODEN: FRMCE8; ISSN: 0014-827X

PUBLISHER: Elsevier Science S.A.

DOCUMENT TYPE: Journal LANGUAGE: English ED Entered STN: 22 Jul 1998

AB (.+-.)-3-Alkoxymethyl-(2-oxabicyclo[3.3.0]octane)-5-yl-methyl-phosphoryl-ethyl-pyridinum [alkyl = Me, CH2=CHCH2, Pr, Me(CH2)5] (I), structurally designed as conformationally restricted platelet activating factor (PAF) antagonists were synthesized in 12-26% overall yield, using Et (.+-.)-3-hydroxymethyl-5-(2- oxabicyclo[3.3.0]octane)carboxylate as key intermediate. The anti-platelet profile of I was evaluated in a PAF-induced aggregation model in rabbit platelet-rich plasma; only I [alkyl = Me(CH2)5] exhibited a modest activity.

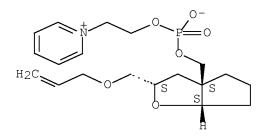
IT 212479-12-4P

(synthesis and platelet activating factor antagonist activity of a new class of bicyclic phospholipids)

RN 212479-12-4 HCAPLUS

CN Pyridinium, 1-[2-[[[(2R,3aR,6aR)-hexahydro-2-[(2-propenyloxy)methyl]-3aH-cyclopenta[b]furan-3a-yl]methoxy]hydroxyphosphinyl]oxy]ethyl]-, inner salt, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



CC 26-3 (Biomolecules and Their Synthetic Analogs)

Section cross-reference(s): 1

IT 197459-82-8P 197459-83-9P 212479-12-4P 212479-14-6P 212479-16-8P 212479-56-6P 212479-58-8P 212479-60-2P

(synthesis and platelet activating factor antagonist activity of a new class of bicyclic phospholipids)

OS.CITING REF COUNT: 7 THERE ARE 7 CAPLUS RECORDS THAT CITE THIS

RECORD (7 CITINGS)

REFERENCE COUNT: 22 THERE ARE 22 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 25 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1997:532277 HCAPLUS Full-text

DOCUMENT NUMBER: 127:166876

ORIGINAL REFERENCE NO.: 127:32228h,32229a

TITLE: Copolymers having phospholipid-like structure and

medical goods coated with them

INVENTOR(S): Shudo, Kenshiro; Matsuyama, Kazuo; Sakaki,

Hidejiro; Kamenosono, Koji; Nakabayashi, Norio;

Ishihara, Kazuhiko

PATENT ASSIGNEE(S): Nippon Oil and Fats Co., Ltd., Japan; Research

Development Corp. of Japan; Nakabayashi, Norio;

Ishihara, Kazuhiko

SOURCE: Jpn. Kokai Tokkyo Koho, 13 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 09183819	Α	19970715	JP 1995-342470	19951228
PRIORITY APPLN. INFO.:			< JP 1995-342470	19951228

ED Entered STN: 20 Aug 1997

Title copolymers have no.-av. mol. wt. 5000-300,000, contain structure units AΒ [CHZCR1[X(Y)mP(O)(O-)O(CH2)2N+R3R4R5]]a[CH2CR9(WSiR6R7R8)]b[R1, R9 = H, Me;R3-R5 = H, C1-6 (hydroxy) hydrocarbyl; R6 = C1-10 alkoxy, C6-14 aryloxy; R7, R8 = C1-10C1-10 (halo) alkoxy, C6-14 (halo) aryloxy, (O- or N-contg.) C1-10 alkyl; W = (CH2)k, CO2(CH2)k, CONH(CH2)k, C6H4(CH2)k, C6H4CH2NH(CH2)k; X = divalent residue; Y = C1-6alkyleneoxy; Z = H, R2O2C; R2 = C1-10 (hydroxy)alkyl; k = 0-4; m = 0-10, a:b = 0(95:5)-(50:50)], and are manufd. by random or alternating polymn. of ZCH:CR1X(Y)mP(O)(O-)O(CH2)2R+R3R4R5 (R1, R3-R5, X, Y, Z, m = same as above) with CH2:CR9WSiR6R7R8 (R6-R9, W = same as above). The copolymers show high durability and biocompatibility. A monomer mixt. comprising 0.6 mol part 2-(methacryloyloxy)ethyl 2'-(trimethylammonio)ethyl phosphate and 0.4 mol part 3-(methacryloyloxypropyl)trimethoxysilane was polymd. at 60.degree. in EtOH using t-Bu peroxypivalate to give a copolymer (Mn 58,000), which was dissolved into EtOH and mixed with H2O and AcOH to prep. a coating. A cover glass was coated with the coating to show protein adsorption 4 ng/cm2.

IT 193684-55-8P

(durable and biocompatible copolymers having phospholipid-like structure for coatings for medical goods)

RN 193684-55-8 HCAPLUS

CN 3,5,8-Trioxa-4-phosphaundec-10-en-1-aminium, 4-hydroxy-N,N,N-trimethyl-, inner salt, 4-oxide, polymer with silicic acid (H4SiO4) ethenyl trimethyl ester (9CI) (CA INDEX NAME)

CM 1

CRN 183544-44-7 CMF C10 H22 N O5 P

CM 2

CRN 19916-97-3 CMF C5 H12 O4 Si

IPCR A61L0033-00 [I,A]; C08F0030-02 [I,A]; C08F0030-08 [I,A]; C08F0230-02
[I,A]; C08F0230-08 [I,A]; C09D0137-00 [I,A]; C09D0143-00 [I,A]

CC 63-8 (Pharmaceuticals)

Section cross-reference(s): 35, 42

IT 193684-52-5P 193684-55-8P 193684-57-0P 193684-58-1P 193684-59-2P

(durable and biocompatible copolymers having phospholipid-like structure for coatings for medical goods)

OS.CITING REF COUNT: 4 THERE ARE 4 CAPLUS RECORDS THAT CITE THIS RECORD (4 CITINGS)

L48 ANSWER 26 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1997:411048 HCAPLUS $\underline{\text{Full-text}}$

DOCUMENT NUMBER: 127:104339

ORIGINAL REFERENCE NO.: 127:19947a,19950a

TITLE: Stimulation of proliferation of V.gamma.2V.delta.2

T cells by alkyl and alkenyl pyrophosphates

INVENTOR(S): Bloom, Barry R.; Tanaka, Yoshimasa; Sano,

Shigetoshi

PATENT ASSIGNEE(S): Albert Einstein College of Medicine of Yeshiva

University, USA

SOURCE: U.S., 25 pp., Cont.-in-part of U.S. Ser. No.

93,528, abandoned.

CODEN: USXXAM

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 5639653	A	19970617	US 1995-390881	19950217
			<	
US 5902793	А	19990511	US 1997-877011	19970616
			<	
PRIORITY APPLN. INFO.:			US 1993-93528	B2 19930719
			<	
			US 1995-390881	A3 19950217

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

Entered STN: 03 Jul 1997

AΒ The proliferation of V.gamma.2V.delta.2 T cells can be stimulated by contacting V.gamma.2V.delta.2 T cells with a compd. selected from the group consisting of a monoalkyl pyrophosphate or an alkenyl pyrophosphate. Thus, 5 .mu.M monoethyl pyrophosphate or 0.3 .mu.M 3-methyl-2-hexenyl pyrophosphate are the minimal ligand concns. required for half-max. proliferative response of .gamma..delta. T cell clone 12G12.

ΙT 6088-04-6, Allyl pyrophosphate

(stimulation of proliferation of V.gamma.2V.delta.2 T cells by

alkyl and alkenyl pyrophosphates)

RN 6088-04-6 HCAPLUS

Diphosphoric acid, P-2-propen-1-yl ester (CA INDEX NAME) CN

INCL 514102000

IPCI C12N0005-02 [ICM, 6]; C12N0005-06 [ICS, 6]; A61K0031-66 [ICS, 6];

C07F0009-02 [ICS,6]

IPCR C07F0009-09 [I,A]; C12N0005-02 [I,A]

NCL 514/102.000; 435/375.000; 435/384.000; 514/106.000; 514/134.000; 558/155.000; 558/156.000

CC 1-7 (Pharmacology)

358-71-4, Isopentenyl pyrophosphate 358-72-5, Dimethylallyl ΙT pyrophosphate 372-97-4, Farnesyl pyrophosphate 763-10-0, Geranyl pyrophosphate 2466-09-3D, Diphosphoric acid, monoalkyl and

monoalkenyl esters 6088-04-6, Allyl pyrophosphate

6699-20-3, Geranylgeranyl pyrophosphate 20680-57-3, Diphosphoric acid, monoethyl ester 22342-44-5, Crotyl pyrophosphate

24753-28-4 52811-47-9, Diphosphoric acid, mono(butyl) ester 56399-35-0, Diphosphoric acid, monomethyl ester 56399-36-1,

Diphosphoric acid, mono(1-methylethyl) ester 104072-24-4,

Diphosphoric acid Mono(Propyl) ester

(stimulation of proliferation of V.gamma.2V.delta.2 T cells by

alkyl and alkenyl pyrophosphates)

11 THERE ARE 11 CAPLUS RECORDS THAT CITE THIS OS.CITING REF COUNT: RECORD (15 CITINGS)

L48 ANSWER 27 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN 1997:94093 HCAPLUS Full-text ACCESSION NUMBER:

DOCUMENT NUMBER: 126:104365

ORIGINAL REFERENCE NO.: 126:20149a,20152a

TITLE: Preparation of substituted liposaccharide analogs

useful in the treatment and prevention of

endotoxemia

INVENTOR(S): Christ, William J.; Rossignol, Daniel P.;

Kobayashi, Seiichi; Kawata, Tsutomu

PATENT ASSIGNEE(S): Eisai Co., Ltd., Japan

PCT Int. Appl., 94 pp. CODEN: PIXXD2 SOURCE:

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 4

PATENT INFORMATION:

PATENT NO.	KIND DATE	APPLICATION NO.	DATE
WO 9639411		WO 1996-US9578	
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US 5681824	A 19971028	US 1995-461677 <	19950605
US 5750664	A 19980512	US 1995-461675 <	19950605
CA 2223140	A1 19961212	CA 1996-2223140 <	19960605
CA 2223140 AU 9663802	C 20080805 A 19961224		19960605
		<	19900003
AU 707779 ZA 9604666	B2 19990722 A 19970311	ZA 1996-4666	19960605
EP 853627	A1 19980722	< EP 1996-923234	19960605
R: AT, BE, CH,	B1 20040121 DE, DK, ES, FR,	< GB, GR, IT, LI, LU, NL,	SE, PT,
IE, FI CN 1192216	A 19980902	CN 1996-195890 <	19960605
CN 1067082 HU 9802662	C 20010613 A2 19990528		19960605
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JP 11506793	T 19990615	JP 1997-501868 <	19960605
JP 4009318 RU 2170738	B2 20071114 C2 20010720		19960605
AT 258185	T 20040215	< AT 1996-923234	19960605
PT 853627	E 20040531	< PT 1996-923234	19960605
ES 2214543	T3 20040916	< ES 1996-923234	19960605
IL 149971	A 20100428		19960605
NO 9705644	A 19980204	< NO 1997-5644 <	19971204

NO 310239	B1	20010611				
JP 2007269812	A	20071018	JP	2007-154127		20070611
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JP 4712001	B2	20110629				
JP 2011121970	A	20110623	JP	2011-18401		20110131
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PRIORITY APPLN. INFO.:			US	1995-461675	A1	19950605
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			JΡ	1997-501868	А3	19960605
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			WO	1996-US9578	M	19960605
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			JP	2007-154127	A3	20070611

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 126:104365

ED Entered STN: 10 Feb 1997

GΙ

ΙT

185955-33-3P

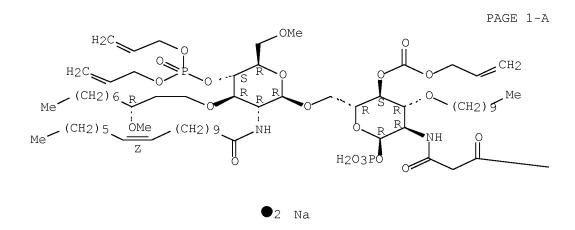
AB Novel substituted liposaccharides in the prophylactic and affirmative treatment of endotoxemia including sepsis, septicemia, and various forms of septic shock and methods of using these agents are provided. Also provided are method of prepg. these agents and intermediates useful therein. Thus, total prepn. of amidodeoxy oligosaccharide I is reported. I inhibited tumor-necrosis factor prodn. in vivo in mice (ED50 = 5 and 10.6 .mu.g/ mouse).

(prepn. of substituted liposaccharide analogs useful in the treatment and prevention of endotoxemia)

RN 185955-33-3 HCAPLUS

CN .alpha.-D-Glucopyranose, 6-O-[4-O-[bis(2-propenyloxy)phosphinyl]-2-deoxy-3-O-[(3R)-3-methoxydecyl]-6-O-methyl-2-[[(11Z)-1-oxo-11-octadecenyl]amino]-.beta.-D-glucopyranosyl]-3-O-decyl-2-deoxy-2-[(1,3-dioxotetradecyl)amino]-, 1-(dihydrogen phosphate) 4-(2-propenyl carbonate), disodium salt (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry as shown.



PAGE 1-B

IPCI C07H0005-04 [ICM,6]; C07H0015-00 [ICS,6]; C07H0017-00 [ICS,6]
IPCR C07H0015-12 [I,A]; A61K0031-70 [I,A]; A61K0031-7028 [I,A]; A61P0031-00
 [I,A]; A61P0031-04 [I,A]; C07H0003-04 [I,A]; C07H0011-00 [I,A];
 C07H0011-04 [I,A]; C07H0013-06 [I,A]; C07H0015-04 [I,A]

CC 33-7 (Carbohydrates)

Section cross-reference(s): 1

IT 41233-29-8P 95548-26-8P 128313-03-1P 128387-27-9P 137766-83-7P 139629-59-7P 139686-99-0P 185954-74-9P 185954-75-0P 185954-76-1P 185954-77-2P 185954-78-3P 185954-79-4P

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185954-80-7P 185954-81-8P
                          185954-82-9P 185954-83-0P
185954-84-1P 185954-85-2P 185954-86-3P 185954-87-4P
185954-88-5P 185954-89-6P 185954-90-9P 185954-91-0P
185954-92-1P 185954-93-2P 185954-94-3P 185954-95-4P
185954-96-5P 185954-97-6P 185955-11-7P 185955-12-8P
185955-13-9P 185955-14-0P 185955-15-1P 185955-16-2P
185955-17-3P 185955-18-4P 185955-19-5P 185955-20-8P
185955-21-9P 185955-22-0P 185955-23-1P 185955-24-2P
185955-25-3P 185955-26-4P 185955-28-6P 185955-29-7P
185955-30-0P 185955-31-1P 185955-32-2P 185955-33-3P
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(prepn. of substituted liposaccharide analogs useful in the treatment and prevention of endotoxemia)

OS.CITING REF COUNT: 6 THERE ARE 6 CAPLUS RECORDS THAT CITE THIS RECORD (10 CITINGS)

REFERENCE COUNT: THERE ARE 2 CITED REFERENCES AVAILABLE FOR 2 THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 28 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1996:713038 HCAPLUS Full-text

125:330468 DOCUMENT NUMBER:

ORIGINAL REFERENCE NO.: 125:61899a,61902a

Low-toxicity aqueous solution of phosphorylcholine TITLE:

group-bearing polymer and its manufacture

INVENTOR(S): Nakabayashi, Nobuo; Ishihara, Kazuhiko; Shuto,

Kenshiro; Matsuyama, Kazuo

PATENT ASSIGNEE(S): Nof Corporation, Japan; Research Development

Corporation of Japan

SOURCE: PCT Int. Appl., 46 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

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	WO	9631	566			A1	_	1996	1010		WO	1996- -1996	JP894 	4		1	9960402
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-		3532				B2		2004				1006	00771	- 0		1	0060400
Ŀ	žΡ	7672	12			AI		1997	0409		EΡ	1996-	90 / /: 	02		1	9960402
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Ţ	JS	6204	324			В1		2001	0320		US	1996-)2		1	9961202
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											JP	1996-		L	i	A 1	9960401
											WO	1996-	JP894	4	Ţ	W 1	9960402

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

ED Entered STN: 05 Dec 1996

AB A process for producing a title soln. useful for skin cares, hair cares, contact lens soilproofing, etc. (no data), comprises polymg. a polymerizable compn. contg. a phosphorylcholine group-bearing monomer of

ZCH:C(R1)X(Y)mP(O)(O-)O(CH2)2N+R2R3R4 [X = divalent org. groups; Y = C1-6 alkylene oxide groups; Z = H; R5OCO, with R5 = C1-10 (hydroxy)alkyl; R1 = H, Me, R1-4 = H, C1-6 (hydroxy)hydrocarbyl; m = 0, 1] in the presence of a nonmetallic polymn. initiator sol. in a water-contg. medium, then purifying the resulting crude aq. soln. by a sepn. membrane. Aq. soln. produced by this process has an impurity content of .1toreq.2000 ppm. Thus, polymg.

2-(methacryloyloxy)ethyl-2'-(trimethylammonio)ethyl phosphate using succinyl peroxide in water gave a polymer which was purified by a dialysis membrane.

IT 183544-45-8P

(low-toxicity aq. soln. of phosphorylcholine group-bearing polymer and manuf.)

RN 183544-45-8 HCAPLUS

CN 3,5,8-Trioxa-4-phosphaundec-10-en-1-aminium, 4-hydroxy-N,N,N-trimethyl-, inner salt, 4-oxide, homopolymer (9CI) (CA INDEX NAME)

CM 1

CRN 183544-44-7 CMF C10 H22 N O5 P

IPCI C08L0043-02 [ICM,6]; C08F0006-06 [ICS,6]; C08F0030-02 [ICS,6]

IPCR C08F0006-00 [I,A]; C08F0006-06 [I,A]; C08F0030-02 [I,A]

CC 37-3 (Plastics Manufacture and Processing)

Section cross-reference(s): 62, 63

IT 67881-99-6P 125275-25-4P 148569-41-9P 150120-18-6P

(low-toxicity aq. soln. of phosphorylcholine group-bearing polymer and manuf.)

OS.CITING REF COUNT: 4 THERE ARE 4 CAPLUS RECORDS THAT CITE THIS

RECORD (9 CITINGS)

REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 29 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1996:115532 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 124:234032

ORIGINAL REFERENCE NO.: 124:43371a,43374a

TITLE: Phosphate multivalent metal salts-modified organopolysiloxanes, their manufacture, and

gelling agents

INVENTOR(S): Ihara, Takeshi; Yano, Shinji; Kita, Katsumi

PATENT ASSIGNEE(S): Kao Corp, Japan

SOURCE: Jpn. Kokai Tokkyo Koho, 13 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 07316169	 А	19951205	JP 1994-112945	19940526
JP 3511401 PRIORITY APPLN. INFO.:	В2	20040329	< JP 1994-112945	19940526
			<	

ED Entered STN: 24 Feb 1996

GΙ

$$-\text{CH}_2\text{CH}_2\text{R}^3 \text{ (OR}^1)_{p0}$$
 PO
 $-\text{(M}^{n+})_{1/n}$
 $\text{R}^4 \text{ (OR}^2)_{q0}$

AB Title siloxanes, useful for gelling agents of silicone oils, cosmetics, medical materials, etc., have .gtoreq.1 Si modified with title salts (I) [R1, R2 = C2-20 linear or branched alkylene; R3 = (hydroxy-substituted) C1-20 linear or branched alkyl; M = alk. earth metal or .gtoreq.2-valent transition metal; p, q = 0-200, n = metal valence no.]. Thus, 10 g phosphate Ca salt and 5.0 g 1,1,1,3,5,5,5-heptamethyltrisiloxane were mixed at 70.degree. for 4h to give a modified organopolysiloxane, which was mixed with octamethylcyclotetrasiloxane to give a gel.

IT 173787-12-7

(manuf. of phosphate metal salt-modified organopolysiloxanes useful for gelling agents of silicone oils and cosmetics and medical materials)

RN 173787-12-7 HCAPLUS

CN Phosphoric acid, monododecyl mono[9-(2-propenyloxy)nonyl] ester, calcium salt (9CI) (CA INDEX NAME)

●1/2 Ca

IPCR C09K0003-00 [I,A]; C07F0009-09 [I,A]; C07F0019-00 [I,A]; C08G0077-38
 [I,A]; C08G0077-395 [I,A]; C08L0083-04 [I,A]; C08L0083-08 [I,A]

CC 37-6 (Plastics Manufacture and Processing)

Section cross-reference(s): 38, 62, 63

IT 1873-88-7, 1,1,1,3,5,5,5-Heptamethyltrisiloxane 17066-04-5 173787-04-7 173787-06-9 173787-08-1 173787-12-7 173787-14-9

(manuf. of phosphate metal salt-modified organopolysiloxanes useful for gelling agents of silicone oils and cosmetics and medical materials)

L48 ANSWER 30 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1995:641004 HCAPLUS Full-text

DOCUMENT NUMBER: 123:199145

ORIGINAL REFERENCE NO.: 123:35573a,35576a

TITLE: Process for preparing phosphinyloxy propanaminium

inner salt derivatives

INVENTOR(S): Prashad, Mahavir; Kapa, Prasad K.

PATENT ASSIGNEE(S): Sandoz Ltd., Switz.

SOURCE: U.S., 20 pp. Continuation of Ser. No. US 93-73407,

filed on 7 Jun 1993, now abando

CODEN: USXXAM

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 5412137	A	19950502	US 1994-197050	19940216
			<	

PRIORITY APPLN. INFO.: US 1993-73407 B1 19930607

<--

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 123:199145

ED Entered STN: 28 Jun 1995

GΙ

Me (CH₂)
$$4^{\circ}$$
 (CH₂) 4° (CH₂) 4° SH N+Me₃

AB The prepn. of the title compds. R1OP(:X1)(X-)OCH(CH2CO2H)CH2N+R2R3R4 where X and X1 are independently O or S; R1 is e.g., alkyl, substituted-alkyl; R2, R3, and R4 are each independently straight or branched chain (C1-4)alkyl, and pharmaceutically acceptable salts, physiol. hydrolyzable esters, and pro-drug

forms thereof, which are useful as hypoglycemic agents (test data given) are described. A representative prepd. compd. is (R)-3-carboxy-N,N,N-trimethyl-2- {[hydroxy(tetradecyloxy)phosphinyl]oxy}-1-propanaminium hydroxide inner salt. 157244-58-1P (prepn. of phosphinyloxy propanaminium inner salts as

hypoglycemics)

RN 157244-58-1 HCAPLUS

ΙT

CN 1-Butanaminium, 2-[[hydroxy(tetradecyloxy)phosphinyl]oxy]-N,N,N-trimethyl-4-oxo-4-(2-propenyloxy)-, inner salt, (R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

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INCL 558146000
IPCI C07F0009-09 [ICM,6]; C07F0009-165 [ICS,6]
IPCR C07F0009-09 [I,A]; C07F0009-113 [I,A]; C07F0009-165 [I,A];
     C07F0009-173 [I,A]
    558/146.000; 558/169.000; 558/170.000
NCL
     29-7 (Organometallic and Organometalloidal Compounds)
CC
     Section cross-reference(s): 1
                                   157244-56-9P
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ΙT
     157244-54-7P
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167685-41-8P 167685-42-9P

(prepn. of phosphinyloxy propanaminium inner salts as

hypoglycemics)

OS.CITING REF COUNT: 3 THERE ARE 3 CAPLUS RECORDS THAT CITE THIS

RECORD (3 CITINGS)

L48 ANSWER 31 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1995:528833 HCAPLUS Full-text

DOCUMENT NUMBER: 123:33387

ORIGINAL REFERENCE NO.: 123:6183a,6186a

TITLE: Preparation of polymerizable phosphorylcholine

derivatives with medical applications

INVENTOR(S): Suzuki, Hiroshi; Kadoma, Yoshihito; Nakabayashi,

Norio; Ishihara, Kazuhiko

PATENT ASSIGNEE(S): Nippon Oils & Fats Co Ltd, Japan SOURCE: Jpn. Kokai Tokkyo Koho, 6 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 07010892	А	19950113	JP 1993-150717	19930622
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JP 3419030	В2	20030623		
PRIORITY APPLN. INFO.:			JP 1993-150717	19930622
			/	

OTHER SOURCE(S): MARPAT 123:33387

ED Entered STN: 06 May 1995

GΙ

$$R^{1}R^{2}R^{3}N^{+}(CH_{2})_{20} - P - O(A)_{m}(0)_{n}COCR^{4} = CH_{2}$$

The title derivs. II (R1-3 = H, C1-3 alkyl; R4 = H, Me; A = linear or branched alkylene; m, n = 0, 1), useful as materials for medical appliances, e.g. catheters, artificial organs and blood vessels, contact lenses, cosmetics, water-absorbing materials, etc., are prepd. by treatment of R1R2R3N+(CH2)2OH Y- (Y- = anion) with X1X2P(O)O(A)m(O)nCOCR4:CH2 (I; X1-2 = halo) in the presence or absence of bases while removing hydrogen halides formed in the reaction followed by hydrolysis of the resulting products by treatment with H2O. This methods give the products without using toxic Me3N with offensive odor. CH2:CMeCO2CH2CH2OH was treated with POC13 and Et3N in CH2C12 at 0.degree. for 4 h to give 95% I (R4 = Me, X1 = X2 = C1, A = CH2CH2, m = n = 1). This was further treated with Me3N+CH2CH2OH C1- and Et3N in CH2C12 at 0.degree. for 2 h, and after removal of Et3N.HC1, treated with H2O at 0.degree. for 2 h to give 65% II (R1 = R2 = R3 = R4 = Me, A = CH2CH2, n = m = 1).

IT 163716-44-7P

(prepn. of (meth) acryloyloxyethyl trialkylammonioethyl phosphates as monomers for medical materials and appliances from (hydroxyethyl) trialkylammonium halides and (meth) acryloyloxyethyl dihalophosphates)

RN 163716-44-7 HCAPLUS

CN 3,5,8-Trioxa-4-phosphaundec-10-en-1-aminium, N,N,N-triethyl-, inner salt, 4-oxide (9CI) (CA INDEX NAME)

IPCI C07F0009-09 [ICM,6]

IPCR C07F0009-09 [I,A]

CC 29-7 (Organometallic and Organometalloidal Compounds)

Section cross-reference(s): 63

IT 163716-43-6P 163716-44-7P

(prepn. of (meth)acryloyloxyethyl trialkylammonioethyl phosphates as monomers for medical materials and appliances from (hydroxyethyl)trialkylammonium halides and (meth)acryloyloxyethyl

dihalophosphates)

L48 ANSWER 32 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1995:526804 HCAPLUS Full-text

DOCUMENT NUMBER: 122:291219

ORIGINAL REFERENCE NO.: 122:53107a,53110a

TITLE: Phosphonooxymethyl or methylthiomethyl ethers of

taxane derivatives as antitumor agents.

INVENTOR(S): Golik, Jerzy; Kadow, John F.; Kaplan, Murray A.;

Li, Wen-Sen; Perrone, Robert K.; Thottathil, John K.; Vyas, Dolatrai; Wittman, Mark D.; Wong, Henry;

Wright, John J.

PATENT ASSIGNEE(S): Bristol-Myers Squibb Co., USA

SOURCE: Eur. Pat. Appl., 124 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 37

PATENT INFORMATION:

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A1	19950222	EP 1994-112803	19940816
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CN	1100771	С	20030205				
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OTHER SOURCE(S): MARPAT 122:291219

ED Entered STN: 05 May 1995

Taxane derivs. T-[OCH2(OCH2)mOP(O)(OH)2]n [T = taxane substituted at C-13 by 3-amino-2-hydroxypropanoyloxy; m = 0-6; n = 1-3] were prepd. from paclitaxel or baccatin III via T'-[OCH2(OCH2)mSMe]n [T' = protected T] for use as antitumor agents. Thus, paclitaxel was converted to its 7-0-methylthiomethyl deriv., treated with dibenzyl phosphate, followed by hydrogenolysis to give 7-0-phosphonooxymethylpaclitaxel. This compd. had an IC50 against HCT-116 human carcinoma of 0.0158 .mu.M. Some of the methylthiomethyl derivs. also have antitumor activity.

IT 1242672-84-9

(Phosphonooxymethyl or methylthiomethyl ethers of taxane derivatives as antitumor agents.)

RN 1242672-84-9 HCAPLUS

CN 2-Furanpropanoic acid, .beta.-[[(1,1-dimethylethoxy)carbonyl]amino]-.alpha.-[[(2-propen-1-yloxy)carbonyl]oxy]-,
(2aR, 4S, 4aS, 6R, 9S, 11S, 12R, 12aR, 12bS)-6, 12b-bis (acetyloxy)-12(benzoyloxy)-2a, 3, 4, 4a, 5, 6, 9, 10, 11, 12, 12a, 12b-dodecahydro-11-hydroxy-4a, 8, 13, 13-tetramethyl-5-oxo-4-[[(phosphonooxy)methoxy]methoxy]-7, 11methano-1H-cyclodeca[3, 4]benz[1, 2-b]oxet-9-yl ester,
(.alpha.R, .beta.R)- (CA INDEX NAME)

Absolute stereochemistry.

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C07F0009-655 [I,A]; C07F0009-6558 [I,A]; H01L0021-48 [I,A]
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     30-20 (Terpenes and Terpenoids)
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       derivatives as antitumor agents.)
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       (Phosphonooxymethyl or methylthiomethyl ethers of taxane
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derivatives as antitumor agents.)

OS.CITING REF COUNT: 24 THERE ARE 24 CAPLUS RECORDS THAT CITE THIS RECORD (26 CITINGS)

L48 ANSWER 33 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1995:297487 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 122:81691

ORIGINAL REFERENCE NO.: 122:15531a,15534a

TITLE: Preparation of phosphonooxymethyl taxane ethers as

neoplasm inhibitors

INVENTOR(S): Golik, Jerzy; Vyas, Dolatrai; Wrigth, John J. Kim;

Wong, Henry; Kadow, John F.; Thotathil, John K.; Li, Wen-Sen; Kaplan, Murray A.; Perrone, Robert K.

PATENT ASSIGNEE(S): Bristol-Myers Squibb Co., USA

SOURCE: Eur. Pat. Appl., 96 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 37

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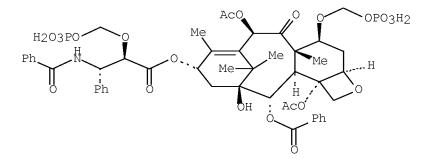
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ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 122:81691

ED Entered STN: 18 Jan 1995

GΙ



AB T[OCH2(OCH2)mOP(O)(OH)2]n [T = a taxane moiety bearing on the C-13 C atom a substituted 3-amino-2-hydroxypropanoyloxy group (sic); m = 0-6; n = 1-3] were prepd. Thus, paclitaxel was converted in 5 steps to 7-O-phosphonooxymethylpaclitaxel triethanolamine salt which gave survival of M109 lung carcinoma-implanted mice 160% that of controls at 24mg/kg i.v. on days 5,6,7,8, and 9 post tumor implant.

IT 1242690-20-5

RN 1242690-20-5 HCAPLUS

CN INDEX NAME NOT YET ASSIGNED

Absolute stereochemistry.

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IPCI C07F0009-655 [ICM,5]; A61K0031-675 [ICS,5]; C07F0009-6558 [ICS,5];
    C07D0305-14 [ICS,5]; C07D0407-12 [ICS,5]; C07F0007-18 [ICS,5]
IPCR A61K0031-665 [I,A]; A61P0035-00 [I,A]; C07D0305-14 [I,A]; C07D0407-12
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    C07F0009-6558 [I,A]; H01L0021-48 [I,A]
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       (Preparation of phosphonooxymethyl taxane ethers as neoplasm
       inhibitors)
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       (Preparation of phosphonooxymethyl taxane ethers as neoplasm
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inhibitors)

OS.CITING REF COUNT: 11 THERE ARE 11 CAPLUS RECORDS THAT CITE THIS RECORD (11 CITINGS)

L48 ANSWER 34 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1994:534463 HCAPLUS Full-text

DOCUMENT NUMBER: 121:134463

ORIGINAL REFERENCE NO.: 121:24325a,24328a

TITLE: Phosphinyloxy propanaminium inner salt derivatives

with hypoglycemic activity

INVENTOR(S): Anderson, Robert Charles; Bebernitz, Gregory R.;

Fraser, James D.; Hughes, Jeffrey W.; Kapa, Prasad K.; Prashad, Mahavir; Smith, Howard C.; Willhauer,

Edwin B.

PATENT ASSIGNEE(S): Sandoz Ltd., Switz.; Sandoz-Patent-G.m.b.H.;

Sandoz-Erfindungen Verwaltungsgesellschaft m.b.H.

SOURCE: Eur. Pat. Appl., 26 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT NO.	KIND	DATE	APPLICATION NO.	
EP 574355			EP 1993-810406 <	
		19970813		
· · · · · · · · · · · · · · · · · · ·	•		GB, GR, IE, IT, LI, LU, HU 1993-1587	
AT 156828	Т	19970815	AT 1993-810406 <	19930607
ES 2105194	Т3	19971016	ES 1993-810406 <	19930607
NO 9302096	A	19931213	NO 1993-2096 <	19930609
AU 9340115	A	19931216	AU 1993-40115 <	19930609
AU 667537	B2	19960328		
IL 105965	A	19970318	IL 1993-105965 <	19930609
CA 2098133	A1	19931212	CA 1993-2098133 <	19930610
JP 06073077	A	19940315	JP 1993-163868 <	19930610
CN 1086217	А	19940504	CN 1993-108729 <	19930610
CN 1039326	С	19980729		
ZA 9304157	A	19940905	ZA 1993-4157 <	19930611
US 5516767	A	19960514	US 1995-373802 <	19950117
PRIORITY APPLN. INFO.:			US 1992-897210 <	A 19920611
			US 1993-72804 <	B1 19930607

US 1994-189856 B1 19940201

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OTHER SOURCE(S): CASREACT 121:134463; MARPAT 121:134463

ED Entered STN: 17 Sep 1994

GΙ

- AB Title compds. ROP(:X1)(X-)OCH(CH2CO2H)CH2N+R1R2R3(X, X1 = 0, S; R = organyl e.g., tetradecyl; R1, R2, R3 = C1-4-alkyl) were prepd. in free acid form or in salt, physiol. hydrolyzable ester or pro-drug form. They can be prepd. by various procedures, e.g. coupling accompanied by oxidn. or thiolation and hydrolysis or thiolysis, or by redn. to amino of a nitro substituent on an arom. ring. They possess hypoglycemic activity and are thus indicated for use in the treatment of diabetes.
- IT 157244-58-1

(prepn. as hypoglycemics)

- RN 157244-58-1 HCAPLUS
- CN 1-Butanaminium, 2-[[hydroxy(tetradecyloxy)phosphinyl]oxy]-N,N,N-trimethyl-4-oxo-4-(2-propenyloxy)-, inner salt, (R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

157244-92-3

IPCR A61K0031-66 [I,A]; A61P0003-08 [I,A]; A61P0003-10 [I,A]; C07C0211-63
 [I,A]; C07F0009-06 [I,A]; C07F0009-09 [I,A]; C07F0009-113 [I,A];
 C07F0009-165 [I,A]

CC 29-7 (Organometallic and Organometalloidal Compounds) Section cross-reference(s): 1

157244-93-4

IT157244-53-6 157244-54-7 157244-55-8 157244-56-9 157244-57-0 157244-58-1 157244-59-2 157244-60-5 157244-61-6 157244-62-7 157244-63-8 157244-64-9 157244-65-0 157244-66-1 157244-67-2 157244-71-8 157244-70-7 157244-68-3 157244-69-4 157244-72-9 157244-73-0 157244-74-1 157244-75-2 157244-76-3 157244-77-4 157244-78-5 157244-79-6 157244-80-9 157244-81-0 157244-82-1 157244-83-2 157244-84-3 157244-85-4 157244-86-5 157244-91-2 157244-87-6 157244-88-7 157244-89-8 157244-90-1

157244-95-6

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157244-97-8 157245-02-8 157245-07-3 157245-12-0 157245-17-5 157245-22-2 157245-27-7 157245-32-4 157245-37-9 157245-42-6	157244-98-9 157245-03-9 157245-08-4 157245-13-1 157245-18-6 157245-23-3 157245-28-8 157245-33-5 157245-38-0 157245-38-0	157244-99-0 157245-04-0 157245-09-5 157245-14-2 157245-19-7 157245-24-4 157245-29-9 157245-34-6 157245-39-1 157245-44-8	157245-00-6 157245-05-1 157245-10-8 157245-15-3 157245-20-0 157245-25-5 157245-30-2 157245-35-7 157245-40-4 157245-45-9	157245-01-7 157245-06-2 157245-11-9 157245-21-1 157245-21-1 157245-31-3 157245-36-8 157245-41-5 157245-46-0
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(prepn. as hypoglycemics)

OS.CITING REF COUNT: 3 THERE ARE 3 CAPLUS RECORDS THAT CITE THIS RECORD (5 CITINGS)

L48 ANSWER 35 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1994:517786 HCAPLUS $\underline{\text{Full-text}}$

DOCUMENT NUMBER: 121:117786

ORIGINAL REFERENCE NO.: 121:21093a,21096a

TITLE: Diester monomer, its polymer, water-containing

soft contact lens, and processing solution for

contact lens

INVENTOR(S): Koinuma, Yasumi; Matsumoto, Takeo; Nakada,

Nobuharu; Nakabayashi, Nobuo; Ishihara, Kazuhiko

PATENT ASSIGNEE(S): NOF Corp., Japan

SOURCE: Eur. Pat. Appl., 26 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 580435	A1	19940126	EP 1993-305795	19930722
EP 580435 R: BE, CH, DE,	B1 FR, GB	19980916 , IT, LI, NL		
JP 06041156	А	19940215	JP 1992-197235 <	19920723
JP 3240695	B2	20011217		
JP 06041157	А	19940215	JP 1992-197238 <	19920723
JP 3240696	В2	20011217		
JP 06043407	А	19940218	JP 1992-197236 <	19920723
JP 3206117	B2	20010904		
JP 06043408	А	19940218	JP 1992-197237 <	19920723
JP 3206118	B2	20010904		
JP 06043400	А	19940218	JP 1992-198633 <	19920724
JP 3227810	В2	20011112		
JP 06043401	A	19940218	JP 1992-198634	19920724

			<		
JP 3227811	В2	20011112			
US 5466853	A	19951114	US 1993-94293		19930719
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PRIORITY APPLN. INFO.:			JP 1992-197235	Α	19920723
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			JP 1992-198634	Α	19920724
			<		

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT OTHER SOURCE(S): MARPAT 121:117786

ED Entered STN: 03 Sep 1994

AB A diester monomer is represented by the formula HR1C:CR3(CH2)mR2 (I), wherein R1, R2, and R3 represent a halogen atom, an alkoxy, an alkenyloxy, a hydroxyalkyloxycarbonyl, or a phosphorylcholine deriv. group. A polymer and a water-contg. contact lens are obtained by polymg. a starting component material including the above-mentioned diester monomer of the formula I. A contact lens processing soln. includes the polymer and a solvent for dissolving the polymer. The polymers provide a soft contact lens with an improved water content and O permeability. For example,

.alpha.-isopropyl-.beta.-[(2'-trimethylammonio)ethyl Et phosphate]itaconate-allyl methacrylate copolymer was prepd. and showed a higher water content and O permeation coeff., compared to those of 2-hydroxyethyl methacrylate-ethylene glycol dimethacrylate copolymer.

IT 156526-61-3P

(prepn. of, for manuf. of soft contact lenses)

RN 156526-61-3 HCAPLUS

CN 3,5,10,15-Tetraoxa-4-phosphaoctadeca-12,17-dien-1-aminium, 4-hydroxy-N,N,N-trimethyl-11,14-dioxo-, inner salt, 4-oxide, (E)-, polymer with 1,2-ethanediyl bis(2-methyl-2-propenoate) and 2-hydroxyethyl 2-propenoate (9CI) (CA INDEX NAME)

CM 1

CRN 156526-60-2 CMF C16 H28 N O8 P

Double bond geometry as shown.

CM 2

CRN 818-61-1 CMF C5 H8 O3

CM 3

CRN 97-90-5 CMF C10 H14 O4

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IPCI C07F0009-10 [ICM,5]; C08F0030-02 [ICS,5]; G02C0007-04 [ICS,5] IPCR C07F0009-09 [I,A]; C08F0030-02 [I,A]; G02B0001-04 [I,A]
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CC 63-7 (Pharmaceuticals)

Section cross-reference(s): 35

IT 156526-57-7P 156526-59-9P 156526-68-0P 156549-95-0P 156549-97-2P 156608-05-8P 156608-07-0P 156608-08-1P 156608-09-2P 156608-11-6P 156608-12-7P 156608-13-8P 156608-14-9P 156608-15-0P 156608-16-1P 156608-17-2P 156608-18-3P 156608-19-4P 156608-20-7P 156608-21-8P 156608-22-9P 156608-24-1P 156608-26-3P 156608-27-4P 156646-89-8P 156646-90-1P 156646-91-2P 156646-93-4P 156978-35-7P

(prepn. of, for manuf. of soft contact lenses)
OS.CITING REF COUNT: 4 THERE ARE 4 CAPLUS RECORDS THAT CITE THIS RECORD (6 CITINGS)

L48 ANSWER 36 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1994:77596 HCAPLUS Full-text

DOCUMENT NUMBER: 120:77596

ORIGINAL REFERENCE NO.: 120:13977a,13980a

TITLE: Preparation and antitumor activity of

anti-endotoxin lipid A analogs

INVENTOR(S): Christ, William J.; Kawata, Tsutomu; Hawkins, Lynn D.; Kobayashi, Seiichi; Asano, Osamu; Rossignol,

Daniel P.

PATENT ASSIGNEE(S): Eisai Co., Ltd., Japan SOURCE: Eur. Pat. Appl., 213 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 536969	A2	19930414	EP 1992-309057	19921005
EP 536969 R: AT, BE, CH,	A3 DE, DK	19940518 , ES, FR, GE	B, GR, IE, IT, LI,	LU, MC, NL,
PT, SE US 5530113	А	19960625	US 1992-935050	19920825
PRIORITY APPLN. INFO.:			< US 1991-776100 <	A 19911011
			US 1992-935050	A 19920825

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 120:77596

ED Entered STN: 19 Feb 1994

GΙ

Title compds. I [R1-R4 = acyl, alkenylcarboxy, alkynylcarboxy; A1, A2 = H, OH, OMe, (CH2)nCO2H, O(CH2)nCO2H, O(CH2)nPO3H2, (CH2)nOPO3H2, n = 0-5; X = H, alkyl, alkenyl, hydroxyalkyl, alkoxyalkyl, (CH2)mOPO3H2, (CH2)mOR5, m = 0-14, R5 = R1-R4; Y = H, OH, halo, OZ(CH2)mMe, Z = bond, CO, CO2], were prepd. as virucides. Thus, compd. II was prepd. and effectively inhibited in vivo LPS-induced prodn. of tumor necrosis factor (TNF) in mice (ED50 = 16.2 .mu.g/mouse).

IT 151663-32-0

(prepn. as intermediate in prepn. of disaccharide lipid A analogs as neoplasm inhibitors) $\,$

RN 151663-32-0 HCAPLUS

CN .alpha.-D-Glucopyranoside, 2-(phosphonooxy)ethyl 2-deoxy-6-0-[2-deoxy-2-[(1,3-dioxotetradecyl)amino]-3-0-[1-oxo-3-[(1-oxo-5-dodecenyl)oxy]decyl]-4-0-phosphono-.beta.-D-glucopyranosyl]-2-[(1,3-dioxotetradecyl)amino]-, 3-(3-hydroxydecanoate) 4-(2-propenyl

carbonate), [3(R), 6[R(Z)]] - (9CI) (CA INDEX NAME)

PAGE 1-B

- CH= CH- (CH₂) 5- Me

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IPCI C07H0013-06 [ICM,5]; C07H0017-04 [ICS,5]; C07H0015-04 [ICS,5];
     C07H0015-06 [ICS,5]; C07H0015-10 [ICS,5]; C07H0015-18 [ICS,5];
     A61K0031-70 [ICS,5]; C07H0013-04 [ICS,5]
IPCR A61K0031-35 [I,A]; A61K0031-351 [I,A]; A61K0031-739 [I,A]; A61P0031-04
     [I,A]; A61P0031-12 [I,A]; C07D0309-10 [I,A]; C07H0003-04 [I,A];
     C07H0005-04 [I,A]; C07H0005-06 [I,A]; C07H0007-02 [I,A]; C07H0011-04
     [I,A]; C07H0013-06 [I,A]; C07H0015-10 [I,A]
CC
     33-4 (Carbohydrates)
     Section cross-reference(s): 1, 6, 34
ΙT
     2430-94-6
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151663-87-5 151663-88-6 151663-89-7 151663-90-0 151663-91-1
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                        152276-43-2
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152276-59-0 152276-61-4 152308-06-0 152308-08-2 152308-10-6
152308-12-8 152308-14-0 152308-16-2 152308-18-4 152308-20-8
152375-80-9 152375-82-1 152376-87-9 152376-89-1 179893-85-7
   (prepn. as intermediate in prepn. of disaccharide lipid A analogs
   as neoplasm inhibitors)
                        RECORD (10 CITINGS)
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THERE ARE 10 CAPLUS RECORDS THAT CITE THIS OS.CITING REF COUNT: 10

L48 ANSWER 37 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1990:514809 HCAPLUS Full-text

DOCUMENT NUMBER: 113:114809

ORIGINAL REFERENCE NO.: 113:19439a,19442a

TITLE: Phenol derivatives as platelet activating factor

inhibitors

Wissner, Allan; Schaub, Robert E.; Sum, Phaik Eng INVENTOR(S):

American Cyanamid Co., USA PATENT ASSIGNEE(S): SOURCE: Eur. Pat. Appl., 88 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 336142	A2	19891011	EP 1989-104429	19890313
			<	
EP 336142	A3	19910424		
R: AT, BE, CH,	DE, ES	FR, GB, GF	R, IT, LI, NL, SE	
NO 8901376	А	19891005	NO 1989-1376	19890331
			<	
AU 8932394	А	19891005	AU 1989-32394	19890403

			<	
AU 626844	В2	19920813		
DK 8901600	А	19891005	DK 1989-1600 <	19890403
FI 8901583	А	19891005	FI 1989-1583	19890403
ZA 8902447	А	19891227	ZA 1989-2447	19890403
JP 02006470	А	19900110	JP 1989-81640 <	19890403
US 5215975	А	19930601	US 1991-763716 <	19910923
US 5234918	А	19930810	US 1991-763714 <	19910923
AU 9228337	А	19930211	AU 1992-28337 <	19921112
US 5411983	А	19950502	US 1993-99037 <	19930728
PRIORITY APPLN. INFO.:			US 1988-177299 <	A 19880404
			US 1989-316721 <	B1 19890303
			US 1991-763714 <	A3 19910923
			US 1991-763716 <	A1 19910923

OTHER SOURCE(S): MARPAT 113:114809

ED Entered STN: 29 Sep 1990

GΙ

Title compds. I [X = alkyl, PhO, Ph, alkoxy, etc.; i = 1-3; j = 1-6; Q = RO, ROCOCH2O, RCO2 (R = H, alkyl, alkenyl); Y = OCO, OCH2, OP(O)(O-)O; R1 = alkyl, alkoxy, halo; n = 1-4; A = 5- to 7-membered arom. heterocyclyl contg. .gtoreq.1N (and which may contain S), (R2)2S+, (R3)3P+; R2 = alkyl; R3 = alkyl, (substituted) Ph] are prepd. Thiazolium phosphate internal salt II (prepd. from PhCHO and glycerol with 7 steps) showed 61.8% inhibition of PAF-induced vascular permeability in guinea pig skin. IT 127164-65-2P

(prepn. and reaction of, in prepn. of platelet activating factor inhibitors)

```
RN 127164-65-2 HCAPLUS
CN Phosphoric acid, mono[3-(bromomethyl)phenyl]
mono[3-(hexadecyloxy)-2-[(2-methyl-2-propen-1-yl)oxy]propyl] ester
(CA INDEX NAME)
```

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IPCI C07F0009-12 [ICM,4]; C07F0009-54 [ICS,4]; C07F0009-65 [ICS,4];
    C07C0149-46 [ICS,4]; A61K0031-66 [ICS,4]; A61K0031-095 [ICS,4]
IPCR A61K0031-095 [I,A]; A61K0031-215 [I,A]; A61K0031-415 [I,A];
    A61K0031-425 [I,A]; A61K0031-426 [I,A]; A61K0031-44 [I,A];
    A61K0031-4418 [I,A]; A61K0031-4425 [I,A]; A61K0031-47 [I,A];
    A61K0031-472 [I,A]; A61K0031-495 [I,A]; A61K0031-50 [I,A];
    A61K0031-505 [I,A]; A61K0031-685 [I,A]; A61P0001-04 [I,A]; A61P0007-00
    [I,A]; A61P0007-02 [I,A]; A61P0009-06 [I,A]; A61P0011-00 [I,A];
    A61P0011-08 [I,A]; A61P0011-16 [I,A]; A61P0017-00 [I,A]; A61P0035-00
    [I,A]; A61P0037-08 [I,A]; C07C0043-174 [I,A]; C07C0043-178 [I,A];
    C07C0059-125 [I,A]; C07C0309-73 [I,A]; C07C0381-00 [I,A]; C07C0381-12
    [I,A]; C07D0213-20 [I,A]; C07D0215-10 [I,A]; C07D0217-10 [I,A];
    C07D0233-60 [I,A]; C07D0233-68 [I,A]; C07D0233-70 [I,A]; C07D0237-08
    [I,A]; C07D0239-20 [I,A]; C07D0239-26 [I,A]; C07D0241-12 [I,A];
    C07D0277-22 [I,A]; C07D0295-04 [I,A]; C07D0521-00 [I,A]; C07F0009-12
    [I,A]; C07F0009-54 [I,A]; C07F0009-58 [I,A]; C07F0009-60 [I,A];
    C07F0009-62 [I,A]; C07F0009-6509 [I,A]; C07F0009-6539 [I,A]
CC
    25-10 (Benzene, Its Derivatives, and Condensed Benzenoid Compounds)
    Section cross-reference(s): 1, 27, 28, 29
               27079-92-1P 54267-06-0P 80350-03-4P 99884-78-3P
IT
    1708-40-3P
    104216-84-4P
                 111841-39-5P
                                111841-45-3P 111841-68-0P
                 127164-51-6P
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    127164-66-3P 127164-67-4P
                                 127164-68-5P 127164-69-6P
    127164-70-9P, 8-Hydroxy-7-methoxy-1-octene 127164-71-0P
    127164-72-1P 127164-73-2P 127164-74-3P
        (prepn. and reaction of, in prepn. of platelet activating factor
       inhibitors)
ΙT
    127164-07-2P 127164-08-3P
                                 127164-09-4P
                                               127164-10-7P
    127164-11-8P 127164-12-9P
                                 127164-13-0P 127164-14-1P
    127164-15-2P 127164-16-3P
                                 127164-17-4P 127164-18-5P
    127164-19-6P 127164-20-9P
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    127164-23-2P 127164-24-3P
                                 127164-25-4P 127164-26-5P
    127164-27-6P 127164-28-7P
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                   127164-36-7P
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    127164-39-0P 127164-40-3P
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    127164-43-6P 127164-44-7P
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    127164-47-0P 127164-48-1P
                                 127164-49-2P 127164-50-5P
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127190-20-9P 127190-21-0P 127190-22-1P 129008-01-1P (prepn. of, as platelet activating factor inhibitor)

OS.CITING REF COUNT: 2 THERE ARE 2 CAPLUS RECORDS THAT CITE THIS

RECORD (2 CITINGS)

L48 ANSWER 38 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1989:57664 HCAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 110:57664

ORIGINAL REFERENCE NO.: 110:9549a,9552a

TITLE: (Pyridinylmethyl) sulfinylbenzimidazole derivatives

as antiulcer agents, their preparation and

formulations containing them

INVENTOR (S): Alminger, Tomas Boerje; Bergman, Rolf Axel;

Bundgaard, Hans; Lindberg, Per Lennart; Sunden,

Gunnel Elisabeth

PATENT ASSIGNEE(S): Aktiebolag Haessle, Swed. SOURCE: PCT Int. Appl., 96 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.				KIND DATE			APPLICATION NO.									
									WO 1987-SE546						9871120	
	:W									FI,	GB, SU,	HU,	JP,	KP,	KR,	LK,
	RW:				CF, TD,			CM,	DE,	FR,	GA,	GB,	IT,	LU,	ML,	MR,
ZA					-			0727	ZA 1987-8263 <				1	9871103		
IL	L 84504			А		1994	1007						1	9871117		
AU	8783	302			А		1988	0616	I	AU 1	-987 >>	8330	2		1	9871120
	6121 2791				B2 A2		1991				-987				1	9871120
											.90/-		02		Τ.	90/1120
ΕP	2791 2791 R:	49					1988									
	2705				A5		1989	0802	Ι		-987 >				1	9871120
ΕP	3326	47			A1		1989	0920	Ε	EP 1		9080			1	9871120
JP	R: 0250									JP 1	NL, -988	5002			1	9871120
HU	5126	9			A2		1990	0428		HU 1		6196			1	9871120
	2048 5107						1992 1992		E	EP 1	-992 >	1088			1	9871120

R: AT, BE, CH, DE, ES, FR, GB, GR, IT, LI, LU, NL, SE

10/596,747

AT	84032	Т	19930115	AT	1987-850362 <		19871120
ES	2052603	Т3	19940716	ES	1987-850362		19871120
RU	2062778	C1	19960627	RU	1987-4614294		19871120
CN	87107309	А	19880824	CN	< 1987-107309		19871121
CN	1023226	С	19931222		<		
PL	157655	B1	19920630	PL	1987-268927		19871121
DK	8803654	А	19880701	DK	1988-3654		19880701
NO	8803229	A	19880916	ИО	1988-3229		19880720
NO	173998	В	19931122		`		
	173998	C	19940302				
	8902454	A	19890519	FI	1989-2454		19890519
US	5215974	A	19930601	US	1991-654394		19910208
LV	10954	В	19960820	LV	1993-1371		19931223
LT	3810	В	19960325	LT	1993-1685		19931228
PRIORITY	APPLN. INFO.:			SE	1986-4998	А	19861121
				SE	1986-5551	А	19861223
				SE	1987-4049	Α	19871016
				ΕP	< 1987-850362	А	19871120
				WO	< 1987-SE546	А	19871120
				US	< 1988-199263	В2	19880518
				US	< 1989-297606	В1	19890113
				US	< 1989-380040 <	В1	19891010

OTHER SOURCE(S): CASREACT 110:57664; MARPAT 110:57664 ED Entered STN: 17 Feb 1989

GΙ

$$R8$$
 $R6$
 $R1$
 $R2$
 $R1$
 $R2$
 $R3$
 $R4$

AΒ The title compds. I [X = S, SO; R1 - R4 = H, C1-8 alkyl, C1-8 alkoxy, alkoxyalkyl,halo, CN, CF3, NO2, etc.; R1R2, R2R3, R3R4 may form (satd. substituted heteroatom-contg. rings); R5 = H, C1-4 alkyl, R6 = H, C1-8 alkyl, C1-8 alkoxy, halo; R7 = H, C1-7 alkyl, C1-7 alkoxy; etc.; R8 = H, C1-8 alkyl, C1-8 alkoxy, halo, etc.; or R6R7 or R7R8 together with the adjacent C atoms in the pyridine ring may form a (satd. O, S, N-contq. ring); Z = OP(:O) (OH) OnR9, OP(:O) (OH) OP(:O) (OH) OnR9, etc.; n = 0 or 1; R9 = H, C1-6 alkyl, etc.], useful as antiulcer agents, were prepd. Pure 1-chloromethyl-6-methoxy-2-[[(4-methoxy-3,5-dimethyl-2pyridinyl)methyl]sulfinyl]-1H-benzimidazole and the mono-triethylammonium salt of phosphoric acid monoethyl ester in CH2Cl2 contq. Et3N were mixed together. solvent was evapd. and the residue was heated at 60.degree. for 5 min. CH2Cl2 was added, distd. off, and the product heated again at 60.degree.. This procedure was repeated 4 times until the reaction was completed. The crude product was purified to give 26% I [X = SO, R1, R2, R4, R5 = H, R3 = R7 = OMe, R6 = R8 = Me, Z = OP(:O)(ONa)OEt] (II). At 1 .mu.mol/kg (administered via the duodenal fistula), II inhibited histamine-induced gastric secretion in dogs by 97%. A syrup contg. sucrose 30, saccharin 0.6, EtOH 5, and II 1, flavoring material 0.05 g, and H2O to 100 mL was prepd.

IT 118381-67-2P

(prepn. of, as antiulcer agent)

RN 118381-67-2 HCAPLUS

CN 1H-Benzimidazole-1-methanol, 6-bromo-2-[[[3,5-dimethyl-4-(2-propenyloxy)-2-pyridinyl]methyl]sulfinyl]-, dihydrogen phosphate (ester), sodium salt (9CI) (CA INDEX NAME)

●x Na

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IPCI C07D0401-12 [ICM, 4]; A61K0031-415 [ICS, 4]
IPCR C07D0401-12 [I,A]; A61K0031-415 [I,A]; A61K0031-4184 [I,A];
     A61K0031-44 [I,A]; A61K0031-4427 [I,A]; A61K0031-675 [I,A];
     A61P0001-04 [I,A]; A61P0043-00 [I,A]; C07D0401-14 [I,A]; C07D0453-02
     [I,A]; C07F0009-6558 [I,A]
CC
     28-9 (Heterocyclic Compounds (More Than One Hetero Atom))
     Section cross-reference(s): 1, 63
     118293-04-2P
ΤТ
                    118293-30-4P
                                   118293-38-2P
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118382-01-7P	118382-02-8P	118382-03-9P	118382-04-0P
118402-37-2P	118402-38-3P	118402-39-4P	118402-40-7P
118402-41-8P			

(prepn. of, as antiulcer agent)

OS.CITING REF COUNT: 15 THERE ARE 15 CAPLUS RECORDS THAT CITE THIS

RECORD (51 CITINGS)

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L48 ANSWER 39 OF 39 HCAPLUS COPYRIGHT 2012 ACS on STN ACCESSION NUMBER: 1981:90388 HCAPLUS Full-text

DOCUMENT NUMBER: 94:90388

ORIGINAL REFERENCE NO.: 94:14641a,14644a
TITLE: Dental cements
PATENT ASSIGNEE(S): Lion Corp., Japan

SOURCE: Jpn. Kokai Tokkyo Koho, 7 pp.

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.		DATE
JP 55139311	Α	19801031	JP 1979-47074		19790417
JP 62057601 PRIORITY APPLN. INFO.:	В	19871202	JP 1979-47074	А	19790417
			<		

ED Entered STN: 12 May 1984

GΙ

Dental cements comprise: (1) reagent A contg. phytic acid [83-86-3] or derivs. (I) (R1, R2, R3, R4 = substituted or nonsubstituted C1-12 alkyl and allyl, or H) and, (2) reagent B contg. alkali metal salts. Reagent A and reagent B are mixed in the presence of H2O for hardening. The cements can be used as fillings. For example, reagent A contg. 50% phytic acid and 40% penta-Na phytate [62989-51-9]

and reagent B contg. silica 30, alumina 18, CaF2 34, AlF3 6 and AlPO4 12 parts were prepd.

IT 76653-86-6

(dental cements contg. alkali metal salts and)

RN 76653-86-6 HCAPLUS

CN D-myo-Inositol, 1,2,4,5,6-pentakis(dihydrogen phosphate) 3-(2-propenyl hydrogen phosphate) (9CI) (CA INDEX NAME)

IPCI A61K0006-08

IPCR A61K0006-06 [I,A]; A61K0006-08 [I,A]

CC 63-7 (Pharmaceuticals)

IT 83-86-3 62989-51-9 **76653-86-6**

(dental cements contg. alkali metal salts and)

=> d his nofile

L34

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T.1
               SEL RN
     FILE 'REGISTRY' ENTERED AT 08:35:53 ON 13 APR 2012
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                OR 107-21-1/BI OR 112-47-0/BI OR 17435-77-7/BI OR
               752234-97-2/BI OR 752234-98-3/BI OR 752234-99-4/BI OR
               752235-00-0/BI OR 855894-56-3/BI OR 855894-57-4/BI OR
               855894-58-5/BI)
L3
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L4
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L6
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L7
         28963 SEA SSS FUL L5
             5 SEA SPE=ON ABB=ON PLU=ON L7 AND L2
L8
             7 SEA SPE=ON ABB=ON PLU=ON L2 NOT L8
L9
               SAV L7 TEMP SAS747/A
               STR
L10
L11
            32 SEA SUB=L7 SSS SAM L10
L12
           797 SEA SUB=L7 SSS FUL L10
L13
           244 SEA SPE=ON ABB=ON PLU=ON L12 AND PMS/CI
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L16
               STR
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L17
L18
         13442 SEA SUB=L7 SSS FUL L16
             2 SEA SPE=ON ABB=ON PLU=ON L18 AND L2
T.19
               SAV L18 TEMP SAS747B/A
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L22
           400 SEA SPE=ON ABB=ON PLU=ON L12
L23
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             1 SEA SPE=ON ABB=ON PLU=ON L22 AND L1
L24
            58 SEA SPE=ON ABB=ON PLU=ON L22 AND PHARM?/SC,SX
L25
L26
            44 SEA SPE=ON ABB=ON PLU=ON L25 AND (1802-2003)/PRY,AY,PY
            37 SEA SPE=ON ABB=ON PLU=ON L22 AND L23
L27
L28
            20 SEA SPE=ON ABB=ON PLU=ON L27 AND (1802-2003)/PRY,AY,PY
L29
            2 SEA SPE=ON ABB=ON PLU=ON L21 AND (1802-2003)/PRY,AY,PY
L30
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               QUE SPE=ON ABB=ON PLU=ON DENTAL? (5A) (ADHES? OR SEAL? OR
L31
               MATERIAL? OR GLUE? OR BOND)
L32
            80 SEA SPE=ON ABB=ON PLU=ON L23 AND L31
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L33
              STR L10
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0 SEA SUB=L12 SSS SAM L33

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			SAV	L35 SAS	747C/A				
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L37		8	SEA	SPE=ON	ABB=ON	PLU=ON	L35		
L38		3	SEA	SPE=ON	ABB=ON	PLU=ON	L37	AND	(1802-2003)/PRY,AY,PY
L39		19	SEA	SPE=ON	ABB=ON	PLU=ON	L30	NOT	L38
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L40			STR	L33					
L41		4	SEA	SUB=L12	SSS SAM	L40			
L42		112	SEA	SUB=L12	SSS FUL	L40			
			SAV	L42 TEM	P SAS747	D/A			
	FILE	'HCAPI	LUS'	ENTERED	AT 11:0	1:34 ON	13 A	PR 20	012
L43		30	SEA	SPE=ON	ABB=ON	PLU=ON	L42		
L44		5	SEA	SPE=ON	ABB=ON	PLU=ON	L43	AND	L26
L45		15	SEA	SPE=ON	ABB=ON	PLU=ON	L43	AND	(1802-2003)/PRY,AY,PY
L46		54	SEA	SPE=ON	ABB=ON	PLU=ON	L45	OR I	L26
L47		48	SEA	SPE=ON	ABB=ON	PLU=ON	L46	NOT	(38 OR L39)
L48		39	SEA	SPE=ON	ABB=ON	PLU=ON	L47	AND	PHARM?/SC,SX
=>									